



Greyhound
CHROMATOGRAPHY
AND ALLIED CHEMICALS

WELLINGTON LABORATORIES

STANDARDS FOR ENVIRONMENTAL
TESTING AND RESEARCH

2016-2018



CERTIFICATE OF REGISTRATION

This is to certify that

Wellington Laboratories

345 Southgate Drive, Guelph, Ontario N1G 3M5 Canada

operates a

Quality Management System

which complies with the requirements of

ISO 9001:2008

for the following scope of registration

The Registration covers the Quality Management System as it applies to the design and provision of reference standards and chemicals for use in environmental analysis and toxicological research.

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Certificate Expiry Date: December 26, 2016

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ISO 9001



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SAI GLOBAL

INFORM. INSPIRE. IMPROVE.

This year marks the 36th anniversary of Wellington Laboratories Inc. (Wellington) and we are proud to present our newest catalogue. Moreover, we are very pleased that we have been able to continue to provide you, our clients, with high quality products, efficient service, and knowledgeable technical support. Thank you for your trust and loyalty, and rest assured, they are well-placed.

In this catalogue you will note that many of the products from our previous catalogue have been retained. This is either due to their popularity, continued scientific relevance, or mandated use in regulations from various jurisdictions. Some products have been discontinued; however, inventory of these products may still be available. Please contact Wellington, or one of our distributors, if you are interested in purchasing withdrawn products and we will do our best to accommodate your request.

You may also note that we have added a number of new products, many of which were announced in our newsletter The Wellington Reporter. You can subscribe to receive these new product announcements by e-mail on our website (www.well-labs.com).

All of our products have been prepared according to detailed procedures that comprise our Quality Management System (QMS) which has been registered to ISO 9001. In addition, Wellington is accredited to ISO Guide 34 for reference material production and ISO/IEC 17025 for our testing and calibration activities. Up-to-date certificates for all of these registrations/accreditations are available upon request.

Finally, as you have or will come to know from our products and research activities, Wellington is a very dynamic company. This is largely due to our staff, who are all highly qualified and have somewhat diverse, yet complimentary backgrounds. Most notably, they are all still excited by the science behind our products.

We trust that our products, service, and overall dedication will give you ...

An Added Measure of Confidence

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As many of you may have noticed, polar bears are often utilized as a Wellington 'sub logo'. They are beautiful animals that are endangered primarily due to human activities. They have also acted as sentinels; contaminants that have been found in their tissues have essentially warned us that certain chemicals are persistent, bioaccumulative, and subject to long range transport. In a way, we are somewhat indebted to them.

The artwork used in this catalogue is dedicated to the polar bear. Each picture is an original created by a local Canadian artist in the medium of their choice. The contributing artists are acknowledged throughout the catalogue. We hope that you enjoy their interpretations and talent.

Front Cover Artwork: Jaqueline Gori - Breslau, Ontario, Canada

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"SUPPORT SOLUTIONS"

Throughout this catalogue, among the products listed are sets of Calibration and Verification Solutions designed for use with a variety of GC/MS applications. These solution sets or kits are denoted by the incorporation of the code CVS into their catalogue numbers, for example **EPA-1613CVS** or **DFP-CVS-B10**.

These calibration solutions are designed to be used with their corresponding "support solutions". These are the solution/mixtures of native or mass-labelled compounds required for sample processing and method validation as determined by the appropriate method.

As an example, the support solutions for **EPA-1613CVS** are:

| | |
|--------------------|--|
| EPA-1613LCS | for sample spiking prior to extraction; |
| EPA-1613CSS | for extract spiking prior to cleanup; |
| EPA-1613ISS | for spiking of the cleaned-up extract prior to analysis; and |
| EPA-1613PAR | a native spiking solution for method performance testing. |

Some sets of calibration solutions in this catalogue (e.g. BDE-CVS-G) were designed by Wellington, with input from clients, and not for a specific method. Support solutions for these are noted where applicable.

GENERAL INFORMATION

WELLINGTON LABORATORIES INC.

Wellington Laboratories Inc. (Wellington) is recognized world-wide as a reputable and trusted source of certified reference standards of organic compounds of environmental concern.

The majority of our products are used for trace analysis or toxicological research and thus their quality is vital. This quality, as well as the success and longevity of the company, is primarily due to our people. Wellington has a well-qualified, dynamic mix of synthetic chemists, analytical chemists and administrative personnel who are committed to customer satisfaction and continual improvement. This is reflected in our ISO registration and accreditations.

THIS CATALOGUE / NEW PRODUCTS

Originally, Wellington's product line was essentially individual standards of polychlorinated dioxins (PCDDs), dibenzofurans (PCDFs) and biphenyls (PCBs). As can be seen from this catalogue, we now offer a very comprehensive selection of native and mass-labelled PCDDs, PCDFs and PCBs as well as ready-to-use calibration and support solutions for:

PCDD/PCDF methods, such as;

- EPA Method 1613B,
- EPA Method 8280,
- EPA Method 8290,
- EPA Method 23,
- JIS Methods K 0311 and K 0312, and,
- European Standard Method 1948-4.

...and, PCB methods, such as;

- EPA Method 1668C,
- Environment Canada Method 1/RM/31,
- JIS Methods K 0311 and K 0312, and,
- European Standard Method 1948-4 (WHO & Marker PCBs).

Throughout this catalogue, you will also find calibration sets for other target analytes such as PBDEs, PFCs, and PAHs as well as multi-point calibration sets for PCDDs, PCDFs and PCBs. We also have combined calibration sets for PCDDs, PCDFs and WHO PCBs, as well as combined sets for the Marker and WHO PCBs (EN 1948-4).

Since our last catalogue, we have added a number of new native and mass-labelled compounds. All of these were announced in a Wellington Reporter, which were posted on our website (www.well-labs.com).

Most notably, added to our product line are:

- Many new native and ¹³C-labelled perfluorinated compounds (PFCs)
- New native and ¹³C-labelled Hexabromocyclododecane mixtures (HBCDs)
- New experimental flame retardants (EFRs)
- New native and mass-labelled organophosphorus compounds
- New native and ¹³C-labelled halogenated carbazoles
- New native telomer acrylates and acetates

...in addition to many other new and creative products.

SYNTHESIS

The chemical standards offered by Wellington are prepared using unambiguous synthetic routes and purified using a battery of methods. All products are purified to a minimum of 98% chemical purity and the isotopic purity of ¹³C-labelled products is required to be > 99%. The structure of all of our compounds is unequivocally confirmed using a variety of techniques including, as appropriate, NMR (400 or 600 MHz), HRGC with LRMS and/or HRMS, UPLC-MS/MS, and SFC/UV/MS/MS.

ACCURACY/TRACEABILITY

All of the solutions listed in this catalogue are prepared in our laboratories using:

- NIST- or NRC-traceable weights for microbalance calibration
- Class A, NIST-traceable volumetric glassware
- Distilled-in-glass or HPLC grade solvents
- Replicate solutions to ensure accuracy and confirm homogeneity

When possible, these solutions are compared to standard reference materials or certified standards from another source. The expanded maximum percent relative uncertainty of solution concentrations is $\pm 5\%$, unless stated otherwise in this catalogue.

VALIDATION/CERTIFICATION

Wellington was the first supplier of PCDDs, PCDFs and WHO PCBs to validate their solution/mixtures using “truly blind” interlaboratory studies. Since 1991, solutions of our PCDD, PCDF and PCB standards have been submitted as part of 30 international round-robins, resulting in over 2000 independent sets of HRGC/HRMS data.

In addition, over the past several years, Wellington has also submitted standards for 13 interlaboratory studies on PBDEs and 17 studies involving PFCs.

Summaries of all of these interlaboratory studies are available on request. However, in all the studies, averages of the data received for all compounds were well within $\pm 10\%$ of the design values.

ANALYSIS/DOCUMENTATION

Each of our products comes with a detailed Certificate of Analysis (CofA) which includes data which the end user should be able to replicate using equivalent instrumentation and conditions. The CofA includes HRGC/LRMS and/or HRGC/HRMS data depending upon the intended use of the product. Those compounds that are not amenable to GC analysis come with LC/MS data.

Additionally, all of our mass-labelled products come with data that clearly shows their isotopic purity. All calibration sets include RRF summaries showing the required linearity. SDSs and handling guidelines are available for all products.

OTHER PRODUCTS/CUSTOM REQUESTS

For products not listed in this catalogue, please visit our website for updates or contact us at info@well-labs.com. Custom solution preparation and synthetic services are also available.

ORDERING, TERMS, WARRANTY & USE

ORDERING INFORMATION

To place an order, please contact the distributor that serves your country. Distributors are listed on pages 10 and 11, as well as on our website.

For Canada, and for other countries where we do not have a distributor, please contact:

Wellington Laboratories Inc.
345 Southgate Drive
Guelph, Ontario CANADA N1G 3M5

Telephone: (519) 822-2436
Toll-free: (800) 578-6985
FAX: (519) 822-2849
E-mail: orders@well-labs.com
Website: www.well-labs.com

When ordering, please provide as much information as possible, including:

- Detailed shipping address and billing address
- Purchase order number, if known
- Catalogue number and description of product
- Quantity and unit size

TERMS & CONDITIONS OF SALE

Prices: A price list for the products listed in this catalogue *is available from your local distributor*. Prices are subject to change without notice.

Payment: Payment terms are net 30 days from date of invoice. Past due invoices will be subject to a 1.5% monthly finance charge.

Note: We may also accept credit card payments.

Shipping & Handling: All shipments will be F.O.B. Guelph, Ontario and will be made using an appropriate courier.

Returns: Please contact Wellington Laboratories Inc. or your local distributor for a return authorization number. No credit or exchange will be approved after 30 days from shipment and without prior authorization.

LIMITED WARRANTY

At the time of shipment, all products are warranted to be free of defects in material and workmanship and to conform to the accompanying technical and purity specifications. Wellington Laboratories Inc. makes no other warranty, expressed or implied, pertaining to the suitability of the product for any specific application. In case of breach of this warranty, the entire liability of Wellington Laboratories Inc. will be limited to the invoice price of the goods. In no case will Wellington Laboratories Inc. be liable for any special, incidental or consequential damages resulting from the use of its products.

ORDERING, TERMS, WARRANTY & USE

INTENDED USE

The products prepared by Wellington Laboratories Inc. are for laboratory use only. They are not for use in humans.

These chemicals should only be used by qualified personnel who are familiar with their potential hazards and are trained in their handling. With all of our products, due care should be exercised to prevent human contact and ingestion.

The absence of a toxicity warning on any of our products must not be interpreted as an indication that there is no possible health hazard.

Safety Data Sheets (SDSs) are supplied upon request.

PACKAGING

For the safety and convenience of our clients, the solutions provided by Wellington Laboratories Inc. are packaged in clear or amber glass flame-sealed ampoules. Crystalline materials are packaged in glass vials with teflon-lined screw caps.

The solution volumes stated in this catalogue are the minimum volumes which will be delivered and should be considered as approximate. To retain the accuracy of the solutions, dilutions should be made using volumetric glassware.

END USE ONLY/NOT FOR RESALE

The reference standards and materials supplied by Wellington Laboratories Inc. are for end use only by the original purchaser and are not to be resold without written authorization from Wellington Laboratories Inc.

DISTRIBUTORS

To determine which distributor serves your country, please visit our website at www.well-labs.com and follow the distributor link under order info.

| DISTRIBUTOR | HEAD OFFICE | CONTACT INFORMATION |
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| WELLINGTON LABORATORIES JAPAN INC. | JAPAN | 1-22-12 Fujimidai Nerima-ku, Tokyo, Japan 177-0034 Phone: +(81) 3-5934-4184 Fax: +(81) 3-5241-4222 Website: www.well-labs.co.jp Email: info@well-labs.co.jp |
| KANTO CHEMICAL CO., INC. | JAPAN | East Muromachi Mitsui BLDG, 2-1, Nihonbashi Muromachi 2-chome, Chuo-ku, Tokyo, Japan 103-0022 Phone: +(81) 3-6214-1090 Fax: +(81) 3-3241-1047 Website: www.kanto.co.jp Email: reag-info@gms.kanto.co.jp |
| WELLINGTON LABORATORIES LLC. | UNITED STATES OF AMERICA | 7208 West 80th Street, Suite 206 Overland Park, KS, USA 66204 Phone: (913) 722-4919 Toll Free: (877) 809-7039 Fax: (913) 722-4669 Website: www.well-labs.com Email: wellington@swbell.net |
| GREYHOUND CHROMATOGRAPHY & ALLIED CHEMICALS | ENGLAND | 6 Kelvin Park, Birkenhead, Merseyside, England CH41 1LT Phone: (+44)-0-151-649-4000 Fax: (+44)-0-151-649-4001 Website: www.greyhoundchrom.com Email: info@greyhoundchrom.com |
| BCP INSTRUMENTS | FRANCE | 12 avenue des Saules 69600 Oullins, France Phone: +33 (0)4 72 49 72 65 Fax: +33 (0)4 72 49 70 45 Website: www.bcp-instruments.com Email: contact@bcp-instruments.com |
| CHEMICAL RESEARCH 2000 S.R.L. | ITALY | Via Santa Margherita di Belice, 16 00133, Rome, Italy Phone: +(39) 06 20630997 Fax: +(39) 06 20685490 Website: www.cr2000.it Email: info@cr2000.it |
| CAMPRO SCIENTIFIC GMBH | GERMANY | Darser Strasse 2A 14167 Berlin, Germany Phone: +49.(0)30.629.01.89.0 Fax: +49.(0)30.629.01.89.89 Website: www.campro.eu Email: info@campro.eu |

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Wellington Laboratories directly via e-mail at: info@well-labs.com

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Judy Riedt
Petersburg, Ontario, Canada

PCDD/PCDF ANALYTICAL METHOD SOLUTIONS

Repeated from our previous catalogues due to their popularity, sets of calibration solutions and support solutions are offered for the following methods:

U.S. EPA Method 1613B
HRGC/HRMS TCDD and TCDF Analysis Solutions
U.S. EPA Method 8280
U.S. EPA Method 8290
U.S. EPA Method 23
European Standard Method EN 1948-4

All of our calibration kits and support solutions are designed and prepared to be used as received according to the procedures set out in their respective methods. They are all accompanied by detailed certificates of analysis which incorporate HRMS data and RRF summaries for the calibration kits.

PCDD/PCDF RESOLUTION TESTING AND WINDOW DEFINING TEST MIXTURES

Also included in this section are the following solution/mixtures of PCDD and PCDF congeners. These are to be used to test and confirm the resolution of the HRGC column being used and to set retention time windows for the PCDD and PCDF congener groups.

| | |
|-----------------|---|
| CS3WT: | EPA-1613CS3 calibration solution combined with PCDD/PCDF window defining congeners and 2378-TCDD resolution testing isomers. |
| 5CWDS: | PCDD/PCDF window defining congener mix. |
| 5TCDD: | 2378-TCDD resolution test mixture. |
| 225TCDF: | 2378-TCDF resolution test mixture. |
| TDTFWD: | Combined PCDD/PCDF window defining and resolution testing mixture for 3 HRGC columns of varying polarity. |



EPA METHOD 1613 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-----------------------|---|-----------------------|
| EPA-1613CVS | EPA Method 1613 Calibration and Verification Solutions CS1-CS5 | 1 kit (5 ampoules) |
| EPA-1613CSL* | CSL Extended Calibration/Low Level | 500 µl |
| EPA-1613CS0.5* | CS0.5 | 500 µl |
| EPA-1613CS1 | CS1 | 500 µl |
| EPA-1613CS2 | CS2 | 500 µl |
| EPA-1613CS3 | CS3 Calibration Verification | 1 ml |
| EPA-1613CS4 | CS4 | 500 µl |
| EPA-1613CS5 | CS5 | 500 µl |

NOTE: 200 µl AMPOULES OF THE CALIBRATION SOLUTIONS ARE ALSO AVAILABLE. PLEASE CONTACT WELLINGTON OR YOUR LOCAL DISTRIBUTOR FOR PRICING INFORMATION.

| | 1613CSL (ng/ml) | 1613CS0.5 (ng/ml) | 1613CS1 (ng/ml) | 1613CS2 (ng/ml) | 1613CS3 (ng/ml) | 1613CS4 (ng/ml) | 1613CS5 (ng/ml) |
|--|--------------------|----------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
| NATIVE PCDDs & PCDFs | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.25 | 0.5 | 2 | 10 | 40 | 200 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| Octachlorodibenzo-p-dioxin | 1.0 | 2.5 | 5.0 | 20 | 100 | 400 | 2000 |
| | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.25 | 0.5 | 2 | 10 | 40 | 200 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| Octachlorodibenzofuran | 1.0 | 2.5 | 5.0 | 20 | 100 | 400 | 2000 |
| LABELLED PCDDs & PCDFs | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 200 | 200 | 200 | 200 | 200 | 200 | 200 |
| | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| CLEANUP STANDARD | | | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.25 | 0.5 | 2 | 10 | 40 | 200 |
| INTERNAL STANDARDS | | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 | 100 | 100 |

* EPA-1613CS0.5 and EPA-1613CSL are not included in the EPA-1613CVS kit and must be ordered separately.

EPA METHOD 1613 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | | | | | Qty/Conc |
|--|---------------------------------------|---------|---------|---------|-----------|----------|
| EPA-1613LCS* | Labelled Compound Stock Solution | | | | | 1.2 ml |
| EPA-1613CSS* | Cleanup Standard Spiking Solution | | | | | 1.2 ml |
| EPA-1613ISS | Internal Standard Spiking Solution | | | | | 1.2 ml |
| EPA-1613PAR* | Precision and Recovery Stock Solution | | | | | 1.2 ml |
| EPA-1613STOCK | EPA Method 1613 Native Stock Solution | | | | | 1.2 ml |
| | | | | | | |
| | 1613LCS | 1613CSS | 1613ISS | 1613PAR | 1613STOCK | |
| NATIVE PCDDs & PCDFs | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | — | — | — | 40 | 400 | |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | — | — | — | 200 | 2000 | |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | — | — | — | 200 | 2000 | |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | — | — | — | 200 | 2000 | |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | — | — | — | 200 | 2000 | |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | — | — | — | 200 | 2000 | |
| Octachlorodibenzo-p-dioxin | — | — | — | 400 | 4000 | |
| | | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | — | — | — | 40 | 400 | |
| 1,2,3,7,8-Pentachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 2,3,4,7,8-Pentachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | — | — | — | 200 | 2000 | |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | — | — | — | 200 | 2000 | |
| Octachlorodibenzofuran | — | — | — | 400 | 4000 | |
| | | | | | | |
| LABELLED PCDDs & PCDFs | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — | — | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — | — | |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — | — | |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — | — | |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — | — | |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 200 | — | — | — | — | |
| | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — | — | |
| | | | | | | |
| CLEANUP STANDARD | | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | — | 40 | — | — | — | |
| | | | | | | |
| INTERNAL STANDARDS | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | — | 200 | — | — | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | — | 200 | — | — | |

* Working solutions are prepared by diluting EPA-1613LCS (in acetone), EPA-1613CSS (in nonane) and EPA-1613PAR (in acetone) 1:50 (v/v)

This solution allows the HRGC/HRMS operator, with one injection, to:

- Set, or confirm, PCDD and PCDF congener group windows
- Test, or confirm, 2,3,7,8-TCDD resolution
- Verify the calibration

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| CS3WT | EPA Method 1613; Calibration and Verification Solution (CS3) combined with Window Defining and 2378-TCDD Resolution Testing Congeners | 500 µl |

| QUANTITATIVE ANALYTES | (ng/ml) | SEMI-QUANTITATIVE ANALYTES | (ng/ml) |
|--|---------|---|---------|
| NATIVE PCDDs & PCDFs: | | WINDOW DEFINERS:* | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 10 | 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 10 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 50 | 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 10 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 50 | 1,2,4,7,9-Pentachlorodibenzo-p-dioxin | 50 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 50 | 1,2,3,8,9-Pentachlorodibenzo-p-dioxin | 50 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 50 | 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin | 50 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin (WD) | 50 | 1,2,3,4,6,7,9-Heptachlorodibenzo-p-dioxin | 50 |
| Octachlorodibenzo-p-dioxin | 100 | | |
| | | 1,3,6,8-Tetrachlorodibenzofuran | 10 |
| 2,3,7,8-Tetrachlorodibenzofuran | 10 | 1,2,8,9-Tetrachlorodibenzofuran | 10 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 50 | 1,3,4,6,8-Pentachlorodibenzofuran | 50 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 50 | 1,2,3,8,9-Pentachlorodibenzofuran | 50 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 50 | 1,2,3,4,6,8-Hexachlorodibenzofuran | 50 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 50 | | |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 50 | 2,3,7,8-TCDD RESOLUTION TESTING ISOMERS: | |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 50 | 1,2,3,4-Tetrachlorodibenzo-p-dioxin | 5 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran (WD) | 50 | 1,2,3,7/1,2,3,8-Tetrachlorodibenzo-p-dioxin mix | 5 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran (WD) | 50 | 1,2,3,9-Tetrachlorodibenzo-p-dioxin | 10 |
| Octachlorodibenzofuran | 100 | | |
| LABELLED PCDDs & PCDFs: | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 200 | | |
| | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | | |
| CLEANUP STANDARD: | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | 10 | | |
| INTERNAL STANDARDS: | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | | |

(WD) - Window Definer

* 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin (last eluting Hexachlorodibenzo-p-dioxin) was not included as it co-elutes with 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin. Use the 1,2,3,4,6,7,9-Heptachlorodibenzo-p-dioxin to set the window.

* 1,2,3,4,8,9-Hexachlorodibenzofuran (last eluting Hexachlorodibenzofuran) was not included as it can interfere with 1,2,3,7,8,9-Hexachlorodibenzofuran. Use the 1,2,3,4,6,7,8-Heptachlorodibenzofuran to set the window.

HRGC/HRMS TCDD/TCDF ANALYSIS SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|---|-----------------------|
| EPA-513CVS | EPA Method 513 Calibration and Verification Solutions CS1-CS6 (6 x 500 µl) | 1 kit (6 ampoules) |
| 513-CS0.25* | CS0.25 Custom Calibration Solution | 500 µl |
| | 513-CS0.25 513-CS1 513-CS2 513-CS3 513-CS4 513-CS5 513-CS6 (ng/ml) (ng/ml) (ng/ml) (ng/ml) (ng/ml) (ng/ml) (ng/ml) | |
| NATIVE TCDD & TCDF | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.25 0.1 0.5 2 10 40 200 | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.25 0.1 0.5 2 10 40 200 | |
| LABELLED TCDD & TCDF | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 100 100 100 100 100 100 | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 100 100 100 100 100 100 | |
| INTERNAL STANDARDS | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 100 100 100 100 100 100 | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 100 100 100 100 100 100 | |
| CLEANUP STANDARD | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | 0.25 0.1 0.5 2 10 40 200 | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 100 100 100 100 100 100 | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 100 100 100 100 100 100 | |
| EPA-513LCSS | EPA Method 513 Labelled TCDD/TCDF Spiking Solution | 1.2 ml |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 100 ng/ml |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 100 ng/ml |
| EPA-513ISS | EPA Method 513 Internal Standard Spiking Solution | 1.2 ml |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 200 ng/ml |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 200 ng/ml |
| EPA-513CSSA | EPA Method 513 Cleanup Standard Spiking Solution | 1.2 ml |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | | 40 ng/ml |
| EPA-513CSSB | EPA Method 513 Alternative Cleanup Standard Spiking Solution | 1.2 ml |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 100 ng/ml |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 100 ng/ml |
| EPA-513PAR | EPA Method 513 Precision and Recovery Solution | 1.2 ml |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | | 40 ng/ml |
| 2,3,7,8-Tetrachlorodibenzofuran | | 40 ng/ml |

* **513-CS0.25** is not included in the EPA-513CVS kit and must be ordered separately.

EPA METHOD 8280 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|---|-----------------------|
| EPA-8280CVS | EPA Method 8280 Calibration and Verification Solutions CC1-CC5 | 1 kit (5 ampoules) |
| EPA-8280CC1 | CC1 | 500 µl |
| EPA-8280CC2 | CC2 | 500 µl |
| EPA-8280CC3 | CC3 Calibration Verification | 1 ml |
| EPA-8280CC4 | CC4 | 500 µl |
| EPA-8280CC5 | CC5 | 500 µl |

NOTE: 200 µl AMPOULES OF THE CALIBRATION SOLUTIONS ARE ALSO AVAILABLE. PLEASE CONTACT WELLINGTON OR YOUR LOCAL DISTRIBUTOR FOR PRICING INFORMATION.

| | 8280CC1 (ng/µl) | 8280CC2 (ng/µl) | 8280CC3 (ng/µl) | 8280CC4 (ng/µl) | 8280CC5 (ng/µl) |
|---|--------------------|--------------------|--------------------|--------------------|--------------------|
| NATIVE PCDDs & PCDFs | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | — | — | 1.25 | — | — |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | — | — | 1.25 | — | — |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| Octachlorodibenzo-p-dioxin | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 2,3,4,7,8-Pentachlorodibenzofuran | — | — | 0.5 | — | — |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | — | — | 1.25 | — | — |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | — | — | 1.25 | — | — |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | — | — | 1.25 | — | — |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | — | — | 1.25 | — | — |
| Octachlorodibenzofuran | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| | | | | | |
| INTERNAL STANDARDS | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| | | | | | |
| CLEANUP STANDARD | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | — | — | 0.25 | — | — |
| | | | | | |
| RECOVERY STANDARDS | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |

EPA METHOD 8280 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc | | | | |
|--|---------------------------------------|----------------|----------------|----------------|----------------|----------------|
| EPA-8280IS | Internal Standard Solution | 1.2 ml | | | | |
| EPA-8280ISB* | Additional Internal Standard Solution | 1.2 ml | | | | |
| EPA-8280CS | Cleanup Standard Solution | 1.2 ml | | | | |
| EPA-8280RS | Recovery Standard Solution | 1.2 ml | | | | |
| EPA-8280MSS | Matrix Spiking Solution | 1.2 ml | | | | |
| | | | | | | |
| | | 8280IS | 8280ISB | 8280CS | 8280RS | 8280MSS |
| | | (ng/μl) | (ng/μl) | (ng/μl) | (ng/μl) | (ng/μl) |
| NATIVE PCDDs & PCDFs | | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | | — | — | — | — | 2.5 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | | — | — | — | — | 6.25 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | | — | — | — | — | — |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | | — | — | — | — | 6.25 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | | — | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | | — | — | — | — | 6.25 |
| Octachlorodibenzo-p-dioxin | | — | — | — | — | 12.5 |
| | | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | | — | — | — | — | 2.5 |
| 1,2,3,7,8-Pentachlorodibenzofuran | | — | — | — | — | 6.25 |
| 2,3,4,7,8-Pentachlorodibenzofuran | | — | — | — | — | — |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | | — | — | — | — | — |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | | — | — | — | — | 6.25 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | | — | — | — | — | — |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | | — | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | | — | — | — | — | 6.25 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | | — | — | — | — | — |
| Octachlorodibenzofuran | | — | — | — | — | 12.5 |
| | | | | | | |
| INTERNAL STANDARDS | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 5.0 | — | — | — | — |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 5.0 | — | — | — | — |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10.0 | — | — | — | — |
| | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 5.0 | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | | 10.0 | — | — | — | — |
| | | | | | | |
| CLEANUP STANDARD | | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | | — | — | 5.0 | — | — |
| | | | | | | |
| RECOVERY STANDARDS | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | — | — | — | 5.0 | — |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | — | — | — | 5.0 | — |
| | | | | | | |
| ADDITIONAL INTERNAL STANDARDS | | | | | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | | — | 5.0 | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | | — | 10.0 | — | — | — |
| | | | | | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | | — | 5.0 | — | — | — |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | | — | 5.0 | — | — | — |
| Octachloro[¹³ C ₁₂]dibenzofuran | | — | 10.0 | — | — | — |

* Not required by US EPA Method 8280

EPA METHOD 8290 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------------|---|--------------------|
| EPA-8290HRCC1-5 | EPA Method 8290 High Resolution Calibration Solutions (HRCC1-HRCC5) | 1 kit (5 ampoules) |
| EPA-8290HRCC0.25* | HRCC0.25 Supplemental Calibration Solution | 500 µl |
| EPA-8290HRCC0.5* | HRCC0.5 Supplemental Calibration Solution | 500 µl |
| EPA-8290HRCC1 | HRCC1 | 500 µl |
| EPA-8290HRCC2 | HRCC2 | 500 µl |
| EPA-8290HRCC3 | HRCC3 Calibration Verification | 1 ml |
| EPA-8290HRCC4 | HRCC4 | 500 µl |
| EPA-8290HRCC5 | HRCC5 | 500 µl |

NOTE: 200 µl AMPOULES OF THE CALIBRATION SOLUTIONS ARE ALSO AVAILABLE. PLEASE CONTACT WELLINGTON OR YOUR LOCAL DISTRIBUTOR FOR PRICING INFORMATION.

| | 8290- HRCC0.25 (ng/ml) | 8290- HRCC0.5 (ng/ml) | 8290- HRCC1 (ng/ml) | 8290- HRCC2 (ng/ml) | 8290- HRCC3 (ng/ml) | 8290- HRCC4 (ng/ml) | 8290- HRCC5 (ng/ml) |
|--|------------------------------|-----------------------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|
| NATIVE PCDDs & PCDFs | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.25 | 0.5 | 1.0 | 2.5 | 10 | 50 | 200 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| Octachlorodibenzo-p-dioxin | 1.25 | 2.5 | 5.0 | 12.5 | 50 | 250 | 1000 |
| NATIVE PCDFs | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.25 | 0.5 | 1.0 | 2.5 | 10 | 50 | 200 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.625 | 1.25 | 2.5 | 6.25 | 25 | 125 | 500 |
| Octachlorodibenzofuran | 1.25 | 2.5 | 5.0 | 12.5 | 50 | 250 | 1000 |
| INTERNAL STANDARDS | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 125 | 125 | 125 | 125 | 125 | 125 | 125 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 125 | 125 | 125 | 125 | 125 | 125 | 125 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 250 | 250 | 250 | 250 | 250 | 250 | 250 |
| INTERNAL PCDFs | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 125 | 125 | 125 | 125 | 125 | 125 | 125 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 125 | 125 | 125 | 125 | 125 | 125 | 125 |
| RECOVERY STANDARDS | | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 125 | 125 | 125 | 125 | 125 | 125 | 125 |

* **EPA-8290HRCC0.25** and **EPA-8290HRCC0.5** are not included in the EPA-8290HRCC1-5 kit and must be ordered separately.

EPA METHOD 8290 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|-------------------------------|----------|
| EPA-8290SFS | Sample Fortification Solution | 1.2 ml |
| EPA-8290RSS | Recovery Standard Solution | 1.2 ml |
| EPA-8290MSS | Matrix Spiking Solution | 1.2 ml |
| EPA-8290STN | Native Stock PCDDs and PCDFs | 1.2 ml |

| | 8290SFS | 8290RSS | 8290MSS | 8290STN |
|--|----------------|----------------|----------------|----------------|
| | (ng/ml) | (ng/ml) | (ng/ml) | (µg/ml) |
| NATIVE PCDDs & PCDFs | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | — | — | 100 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | — | — | 250 | 2.5 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | — | — | 250 | 2.5 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | — | — | 250 | 2.5 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | — | — | 250 | 2.5 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | — | — | 250 | 2.5 |
| Octachlorodibenzo-p-dioxin | — | — | 500 | 5.0 |
| | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | — | — | 100 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzofuran | — | — | 250 | 2.5 |
| 2,3,4,7,8-Pentachlorodibenzofuran | — | — | 250 | 2.5 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | — | — | 250 | 2.5 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | — | — | 250 | 2.5 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | — | — | 250 | 2.5 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | — | — | 250 | 2.5 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | — | — | 250 | 2.5 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | — | — | 250 | 2.5 |
| Octachlorodibenzofuran | — | — | 500 | 5.0 |
| | | | | |
| INTERNAL STANDARDS | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | — | — | — |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 250 | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 250 | — | — | — |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 500 | — | — | — |
| | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | — | — | — |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 250 | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 250 | — | — | — |
| | | | | |
| RECOVERY STANDARDS | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | 500 | — | — |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | 500 | — | — |

EPA METHOD 23 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|-----------------------|
| EPA-23CS1-5 | EPA Method 23 HRGC/HRMS Calibration Solutions CS1-CS5 | 1 kit (5 ampoules) |
| EPA-23CS1 | CS1 | 500 µl |
| EPA-23CS2 | CS2 | 500 µl |
| EPA-23CS3 | CS3 Calibration Verification | 1 ml |
| EPA-23CS4 | CS4 | 500 µl |
| EPA-23CS5 | CS5 | 500 µl |

NOTE: 200 µl AMPOULES OF THE CALIBRATION SOLUTIONS ARE ALSO AVAILABLE. PLEASE CONTACT WELLINGTON OR YOUR LOCAL DISTRIBUTOR FOR PRICING INFORMATION.

| | 23CS1 (ng/ml) | 23CS2 (ng/ml) | 23CS3 (ng/ml) | 23CS4 (ng/ml) | 23CS5 (ng/ml) |
|--|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| NATIVE PCDDs & PCDFs | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.5 | 1 | 5 | 50 | 100 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| Octachlorodibenzo-p-dioxin | 5.0 | 10 | 50 | 500 | 1000 |
| | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.5 | 1 | 5 | 50 | 100 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| Octachlorodibenzofuran | 5.0 | 10 | 50 | 500 | 1000 |
| | | | | | |
| INTERNAL STANDARDS | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 200 | 200 | 200 | 200 | 200 |
| | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 100 | 100 | 100 | 100 | 100 |
| | | | | | |
| SURROGATE STANDARDS | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | 0.5 | 1 | 5 | 50 | 100 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 2.5 | 5 | 25 | 250 | 500 |
| | | | | | |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| | | | | | |
| ALTERNATIVE STANDARD | | | | | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 2.5 | 5 | 25 | 250 | 500 |
| | | | | | |
| RECOVERY STANDARDS | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 100 | 100 | 100 | 100 | 100 |

EPA METHOD 23 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|-----------------------------------|----------|
| EPA-23IS | Internal Standard Solution | 1.2 ml |
| EPA-23ISS | Internal Standard Stock Solution | 1.2 ml |
| EPA-23SS | Surrogate Standard Solution | 1.2 ml |
| EPA-23SSS | Surrogate Standard Stock Solution | 1.2 ml |
| EPA-23RS | Recovery Standard Solution | 1.2 ml |
| EPA-23AS | Alternative Standard | 1.2 ml |

| | 23IS (ng/ml) | 23ISS (µg/ml) | 23SS (ng/ml) | 23SSS (µg/ml) | 23RS (ng/ml) | 23AS (ng/ml) |
|---|------------------------|-------------------------|------------------------|-------------------------|------------------------|------------------------|
| INTERNAL STANDARDS | | | | | | |
| 2,3,7,8-Tetrachloro ¹³ C ₁₂]dibenzo-p-dioxin | 100 | 1.0 | — | — | — | — |
| 1,2,3,7,8-Pentachloro ¹³ C ₁₂]dibenzo-p-dioxin | 100 | 1.0 | — | — | — | — |
| 1,2,3,6,7,8-Hexachloro ¹³ C ₁₂]dibenzo-p-dioxin | 100 | 1.0 | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro ¹³ C ₁₂]dibenzo-p-dioxin | 100 | 1.0 | — | — | — | — |
| Octachloro ¹³ C ₁₂]dibenzo-p-dioxin | 200 | 2.0 | — | — | — | — |
| | | | | | | |
| 2,3,7,8-Tetrachloro ¹³ C ₁₂]dibenzofuran | 100 | 1.0 | — | — | — | — |
| 1,2,3,7,8-Pentachloro ¹³ C ₁₂]dibenzofuran | 100 | 1.0 | — | — | — | — |
| 1,2,3,6,7,8-Hexachloro ¹³ C ₁₂]dibenzofuran | 100 | 1.0 | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro ¹³ C ₁₂]dibenzofuran | 100 | 1.0 | — | — | — | — |
| SURROGATE STANDARDS | | | | | | |
| 2,3,7,8-[³⁷ Cl ₄]-Tetrachlorodibenzo-p-dioxin | — | — | 100 | 1.0 | — | — |
| 1,2,3,4,7,8-Hexachloro ¹³ C ₁₂]dibenzo-p-dioxin | — | — | 100 | 1.0 | — | — |
| | | | | | | |
| 2,3,4,7,8-Pentachloro ¹³ C ₁₂]dibenzofuran | — | — | 100 | 1.0 | — | — |
| 1,2,3,4,7,8-Hexachloro ¹³ C ₁₂]dibenzofuran | — | — | 100 | 1.0 | — | — |
| 1,2,3,4,7,8,9-Heptachloro ¹³ C ₁₂]dibenzofuran | — | — | 100 | 1.0 | — | — |
| ALTERNATIVE STANDARD | | | | | | |
| 1,2,3,7,8,9-Hexachloro ¹³ C ₁₂]dibenzofuran | — | — | — | — | — | 250 |
| RECOVERY STANDARDS | | | | | | |
| 1,2,3,4-Tetrachloro ¹³ C ₁₂]dibenzo-p-dioxin | — | — | — | — | 500 | — |
| 1,2,3,7,8,9-Hexachloro ¹³ C ₁₂]dibenzo-p-dioxin | — | — | — | — | 500 | — |

EUROPEAN METHOD EN-1948 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|---|-----------------------|
| EN-1948CVS | European Method EN-1948 Calibration and Verification Solutions CS1-CS6 | 1 kit (6 ampoules) |
| EN-1948CSL* | CSL Extended Calibration/Low Level | 500 µl |
| EN-1948CS1 | CS1 | 500 µl |
| EN-1948CS2 | CS2 | 500 µl |
| EN-1948CS3 | CS3 | 500 µl |
| EN-1948CS4 | CS4 | 500 µl |
| EN-1948CS5 | CS5 | 500 µl |
| EN-1948CS6 | CS6 | 500 µl |

NOTE: 200 µl AMPOULES OF THE CALIBRATION SOLUTIONS ARE ALSO AVAILABLE. PLEASE CONTACT WELLINGTON OR YOUR LOCAL DISTRIBUTOR FOR PRICING INFORMATION.

| | 1948CSL (pg/µl) | 1948CS1 (pg/µl) | 1948CS2 (pg/µl) | 1948CS3 (pg/µl) | 1948CS4 (pg/µl) | 1948CS5 (pg/µl) | 1948CS6 (pg/µl) |
|--|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
| NATIVE PCDDs & PCDFs | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.04 | 0.2 | 0.8 | 4 | 16 | 80 | 320 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.16 | 0.8 | 3.2 | 16 | 64 | 320 | 1280 |
| Octachlorodibenzo-p-dioxin | 0.16 | 0.8 | 3.2 | 16 | 64 | 320 | 1280 |
| | | | | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.04 | 0.2 | 0.8 | 4 | 16 | 80 | 320 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.08 | 0.4 | 1.6 | 8 | 32 | 160 | 640 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.16 | 0.8 | 3.2 | 16 | 64 | 320 | 1280 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.16 | 0.8 | 3.2 | 16 | 64 | 320 | 1280 |
| Octachlorodibenzofuran | 0.16 | 0.8 | 3.2 | 16 | 64 | 320 | 1280 |
| | | | | | | | |
| SAMPLING STANDARDS | | | | | | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 32 | 32 | 32 | 32 | 32 | 32 | 32 |
| | | | | | | | |
| EXTRACTION STANDARDS | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 32 | 32 | 32 | 32 | 32 | 32 | 32 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 32 | 32 | 32 | 32 | 32 | 32 | 32 |
| | | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 32 | 32 | 32 | 32 | 32 | 32 | 32 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 32 | 32 | 32 | 32 | 32 | 32 | 32 |
| | | | | | | | |
| SYRINGE STANDARDS | | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 16 | 16 | 16 | 16 | 16 | 16 | 16 |

* **EN-1948CSL** is not included in the EN-1948CVS kit and must be ordered separately.

EUROPEAN METHOD EN-1948 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------|------------------------------|----------|
| EN-1948ES | Extraction Standard Solution | 1.2 ml |
| EN-1948IS | Syringe Standard Solution | 1.2 ml |
| EN-1948SS | Sampling Standard Solution | 1.2 ml |
| EN-1948STK | PCDD/PCDF Solution/Mixture | 1.2 ml |

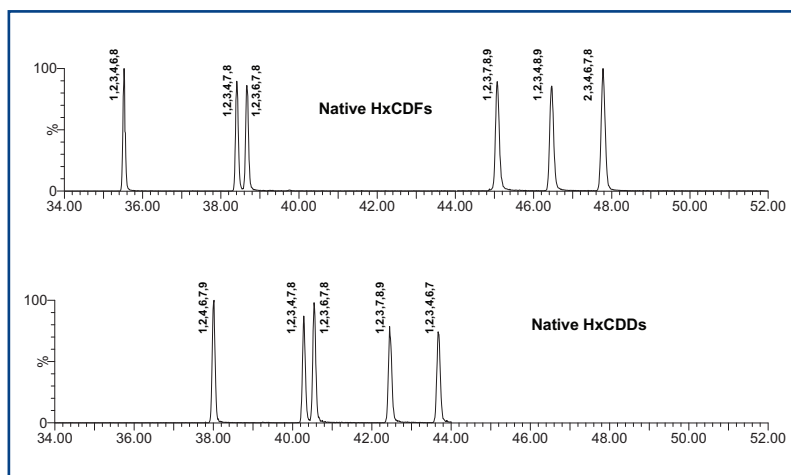
| | 1948ES (pg/μl) | 1948IS (pg/μl) | 1948SS (pg/μl) | 1948STK (μg/ml) |
|--|-------------------|-------------------|-------------------|--------------------|
| NATIVE PCDDs & PCDFs | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | — | — | — | 0.50 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | — | — | — | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | — | — | — | 1.0 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | — | — | — | 1.0 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | — | — | — | 1.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | — | — | — | 2.0 |
| Octachlorodibenzo-p-dioxin | — | — | — | 2.0 |
| | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | — | — | — | 0.50 |
| 1,2,3,7,8-Pentachlorodibenzofuran | — | — | — | 1.0 |
| 2,3,4,7,8-Pentachlorodibenzofuran | — | — | — | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | — | — | — | 1.0 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | — | — | — | 1.0 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | — | — | — | 1.0 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | — | — | — | 1.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | — | — | — | 2.0 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | — | — | — | 2.0 |
| Octachlorodibenzofuran | — | — | — | 2.0 |
| | | | | |
| SAMPLING STANDARDS | | | | |
| 1,2,3,7,8-Pentachloro ^[13C₁₂] dibenzofuran | — | — | 200 | — |
| 1,2,3,7,8,9-Hexachloro ^[13C₁₂] dibenzofuran | — | — | 200 | — |
| 1,2,3,4,7,8,9-Heptachloro ^[13C₁₂] dibenzofuran | — | — | 400 | — |
| | | | | |
| EXTRACTION STANDARDS | | | | |
| 2,3,7,8-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | 200 | — | — | — |
| 1,2,3,7,8-Pentachloro ^[13C₁₂] dibenzo-p-dioxin | 200 | — | — | — |
| 1,2,3,4,7,8-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | 200 | — | — | — |
| 1,2,3,6,7,8-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | 200 | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro ^[13C₁₂] dibenzo-p-dioxin | 400 | — | — | — |
| Octachloro ^[13C₁₂] dibenzo-p-dioxin | 400 | — | — | — |
| | | | | |
| 2,3,7,8-Tetrachloro ^[13C₁₂] dibenzofuran | 200 | — | — | — |
| 2,3,4,7,8-Pentachloro ^[13C₁₂] dibenzofuran | 200 | — | — | — |
| 1,2,3,4,7,8-Hexachloro ^[13C₁₂] dibenzofuran | 200 | — | — | — |
| 1,2,3,6,7,8-Hexachloro ^[13C₁₂] dibenzofuran | 200 | — | — | — |
| 2,3,4,6,7,8-Hexachloro ^[13C₁₂] dibenzofuran | 200 | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro ^[13C₁₂] dibenzofuran | 400 | — | — | — |
| Octachloro ^[13C₁₂] dibenzofuran | 400 | — | — | — |
| | | | | |
| SYRINGE STANDARDS | | | | |
| 1,2,3,4-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | — | 800 | — | — |
| 1,2,3,7,8,9-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | — | 800 | — | — |

CAPILLARY COLUMN PERFORMANCE TEST MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------|
| 5CWDS | Window Defining Mixture for DB-5, BP5, HP-2, Rtx-5, SPB-5, or Equivalent Capillary Columns | 1.2 ml |
| | 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,4,7,9-Pentachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,8,9-Pentachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,4,6,7,9-Heptachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | Octachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,3,6,8-Tetrachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,8,9-Tetrachlorodibenzofuran | 1.0 µg/ml |
| | 1,3,4,6,8-Pentachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,3,8,9-Pentachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,3,4,6,8-Hexachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,3,4,8,9-Hexachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 1.0 µg/ml |
| | Octachlorodibenzofuran | 1.0 µg/ml |
| 5TCDD | 2378-TCDD Isomer Specificity Test Mixture for DB-5, BP5, HP-2, Rtx-5, SPB-5, or Equivalent Columns | 1.2 ml |
| | 1,2,3,4-Tetrachlorodibenzo-p-dioxin | 0.5 µg/ml |
| | 1,2,3,7 and 1,2,3,8-Tetrachlorodibenzo-p-dioxin mix | 0.5 µg/ml |
| | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,9-Tetrachlorodibenzo-p-dioxin | 1.0 µg/ml |
| | 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 µg/ml |
| | 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 0.5 µg/ml |
| 225TCDF | 2378-TCDF Isomer Specificity Test Mixture for DB-225, BP225, HP-225, Rtx-225, SPB-225, or Equivalent Columns | 1.2 ml |
| | 1,3,6,8-Tetrachlorodibenzofuran | 1.0 µg/ml |
| | 2,3,4,7-Tetrachlorodibenzofuran | 1.0 µg/ml |
| | 2,3,7,8-Tetrachlorodibenzofuran | 2.0 µg/ml |
| | 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 0.5 µg/ml |
| | 1,2,3,9-Tetrachlorodibenzofuran | 1.0 µg/ml |
| | 1,2,8,9-Tetrachlorodibenzofuran | 1.0 µg/ml |

CAPILLARY COLUMN PERFORMANCE TEST MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|---|-----------|
| STDWD | Combined Window Defining/TCDD Resolution Testing Mixture for DB-5, BP5, HP-2, Rtx-5, SPB-5, or Equivalent Columns | 1.2 ml |
| WINDOW DEFINERS | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,4,7,9-Pentachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,3,8,9-Pentachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,3,4,6,7,9-Heptachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | | 100 ng/ml |
| Octachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,3,6,8-Tetrachlorodibenzofuran | | 100 ng/ml |
| 1,2,8,9-Tetrachlorodibenzofuran | | 100 ng/ml |
| 1,3,4,6,8-Pentachlorodibenzofuran | | 100 ng/ml |
| 1,2,3,8,9-Pentachlorodibenzofuran | | 100 ng/ml |
| 1,2,3,4,8,9-Hexachlorodibenzofuran | | 100 ng/ml |
| 1,2,3,4,6,8-Hexachlorodibenzofuran | | 100 ng/ml |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | | 100 ng/ml |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | | 100 ng/ml |
| Octachlorodibenzofuran | | 100 ng/ml |
| 2378-TCDD RESOLUTION TESTING ISOMERS | | |
| 1,2,3,4-Tetrachlorodibenzo-p-dioxin | | 50 ng/ml |
| 1,2,3,7 and 1,2,3,8-Tetrachlorodibenzo-p-dioxin mix | | 50 ng/ml |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | | 100 ng/ml |
| 1,2,3,9-Tetrachlorodibenzo-p-dioxin | | 100 ng/ml |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 50 ng/ml |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 50 ng/ml |
| OTHERS | | |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 50 ng/ml |



GC/MS DATA: Native HxCDFs and HxCDDs on a 60m SP-2331 Column.

CAPILLARY COLUMN PERFORMANCE TEST MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------|----------|
|------------------|---------------------------|----------|

| | | |
|---------------|--|--------|
| TDTFWD | Combined Window Defining and Resolution Testing Mixture for 3 Capillary Columns. | 1.2 ml |
|---------------|--|--------|

Concentrations for each compound are listed in brackets (ng/ml ± 20%)^a

WINDOW DEFINING STANDARDS

| 5/2 ^b | | 2331 ^c /225 ^d | |
|--------------------|--------------------|-------------------------------------|--------------------|
| FIRST | LAST | FIRST | LAST |
| 1368-TCDD (50) | 1289-TCDD (50) | 1368-TCDD (50) | 1289-TCDD (50) |
| 12479-PeCDD (50) | 12389-PeCDD (50) | 12479-PeCDD (50) | 12389-PeCDD (50) |
| 124679-HxCDD (50) | 123467-HxCDD (50) | 124679-HxCDD (50) | 123467-HxCDD (50) |
| 1234679-HpCDD (50) | 1234678-HpCDD (50) | 1234679-HpCDD (50) | 1234678-HpCDD (50) |
| | OCDD (50) | | OCDD (50) |
| 1368-TCDF (100) | 1289-TCDF (100) | 1368-TCDF (100) | 1289-TCDF (100) |
| 13468-PeCDF (50) | 12389-PeCDF (50) | 13468-PeCDF (50) | 23467-PeCDF (50) |
| 123468-HxCDF (50) | 123489-HxCDF (50) | 123468-HxCDF (50) | 234678-HxCDF (50) |
| 1234678-HpCDF (50) | 1234789-HpCDF (50) | 1234678-HpCDF (50) | 1234789-HpCDF (50) |
| | OCDF (50) | | OCDF (50) |

RESOLUTION TESTING MIXTURES

| | 5/2 ^b | 2331 ^c | 225 ^d |
|------------------|--|--|--|
| 2378-TCDD | 1234-TCDD (25) 1237 and 1238-TCDD (25) ^e 2378-TCDD (50) 1239-TCDD (50) | 1478-TCDD (25) 2378-TCDD (50) 1237 and 1238-TCDD (25) ^e 1234-TCDD (25) | 1478-TCDD (25) 2378-TCDD (50) 1237 and 1238-TCDD (25) ^e 1234-TCDD (25) |
| 2378-TCDF | 2347-TCDF/ 2348-TCDF/ 2378-TCDF (not resolved) | 1269-TCDF (50) 2378-TCDF (100) 2348-TCDF (50) | 2347-TCDF (50) 2378-TCDF (100) 1239-TCDF (50) |

OTHER PCDDs AND PCDFs INCLUDED

| | | | |
|------------------|-------------------|-------------------|--|
| 12378-PeCDD (50) | 123478-HxCDD (50) | 123478-HxCDF (50) | ¹³ C ₁₂ -1234-TCDD (25) |
| 12378-PeCDF (50) | 123678-HxCDD (50) | 123678-HxCDF (50) | ¹³ C ₁₂ -2378-TCDD (25) |
| 23478-PeCDF (50) | 123789-HxCDD (50) | 123789-HxCDF (50) | ¹³ C ₁₂ -2378-TCDF (25) |
| | | | ¹³ C ₁₂ -123789-HxCDD (50) |

- a Maximum percent relative combined uncertainty of weights and volumes
 b 5/2 - DB-5, BP5, HP-2, Rtx-5, SPB-5 or equivalent capillary column
 c 2331 - SP-2331, Rtx-2330 or equivalent capillary column
 d 225 - DB-225, BP225, HP-225, Rtx-225, SPB-225 or equivalent capillary column
 e Total concentration of both isomers

INDIVIDUAL PCDD & PCDF CONGENERES: NATIVE AND MASS-LABELLED

Wellington has prepared and offers a large selection of individual native and mass-labelled chlorinated dibenzo-p-dioxins (PCDDs) and chlorinated dibenzofurans (PCDFs).

All of these compounds have been synthesized using single product, unambiguous routes and purified using a variety of techniques. Prior to release, their structures and chemical and isotopic purities are confirmed using a number of methods and this data is summarized in the Certificates of Analysis (CofAs).

Accurate solutions of these compounds were prepared as described in the introduction to this catalogue and the maximum percent relative uncertainty associated with their concentrations is $\pm 5\%$. In addition, the concentrations of the 2,3,7,8-substituted PCDDs and PCDFs have been continually certified since 1991 through a large number of interlaboratory studies and are thus traceable to these studies.

The following product groups are presented in this section:

Native Chlorinated Dibenzo-p-dioxins (PCDDs)

Native Chlorinated Dibenzofurans (PCDFs)

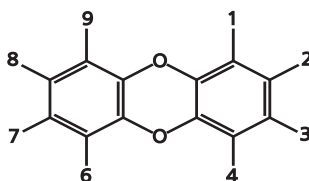
Mass-labelled PCDDs

Mass-labelled PCDFs



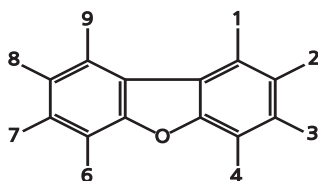
NATIVE CHLORINATED DIBENZO-p-DIOXINS (PCDDs)

| Catalogue Number | Product (toluene or nonane solution) | Qty/Conc |
|----------------------|---|-----------------|
| DD-0-S | Dibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1-S | 1-Chlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-2-S | 2-Chlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-23-S | 2,3-Dichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-27-S | 2,7-Dichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-28-S | 2,8-Dichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123-S | 1,2,3-Trichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-124-S | 1,2,4-Trichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-237-S | 2,3,7-Trichlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1234-S | 1,2,3,4-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1247/8-S | 1,2,4,7/1,2,4,8-Tetrachlorodibenzo-p-dioxin mix | 1.2 ml 50 µg/ml |
| DD-1278-S | 1,2,7,8-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1289-S | 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1368-S | 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1378-S | 1,3,7,8-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1379-S | 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1478-S | 1,4,7,8-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-2378-S | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-12378-S | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-12478-S | 1,2,4,7,8-Pentachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123467-S | 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123468-S | 1,2,3,4,6,8-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123478-S | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123678-S | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-123789-S | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-124679-S | 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-1234678-S | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| DD-12346789-S | Octachlorodibenzo-p-dioxin | 1.2 ml 50 µg/ml |



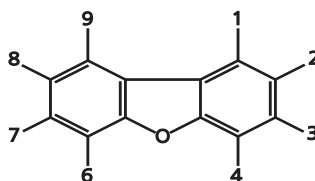
NATIVE CHLORINATED DIBENZOFURANS (PCDFs)

| Catalogue Number | Product (toluene or nonane solution) | Qty/Conc |
|------------------|--------------------------------------|-----------------|
| DF-0-S | Dibenzofuran | 1.2 ml 50 µg/ml |
| DF-2-S | 2-Chlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-4-S | 4-Chlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-23-S | 2,3-Dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-24-S | 2,4-Dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-26-S | 2,6-Dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-27-S | 2,7-Dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-28-S | 2,8-Dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-136-S | 1,3,6-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-138-S | 1,3,8-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-146-S | 1,4,6-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-147-S | 1,4,7-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-149-S | 1,4,9-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-234-S | 2,3,4-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-236-S | 2,3,6-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-238-S | 2,3,8-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-246-S | 2,4,6-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-247-S | 2,4,7-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-248-S | 2,4,8-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-267-S | 2,6,7-Trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1236-S | 1,2,3,6-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1238-S | 1,2,3,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1239-S | 1,2,3,9-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1246-S | 1,2,4,6-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1247-S | 1,2,4,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1248-S | 1,2,4,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1267-S | 1,2,6,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1278-S | 1,2,7,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1279-S | 1,2,7,9-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1347-S | 1,3,4,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |



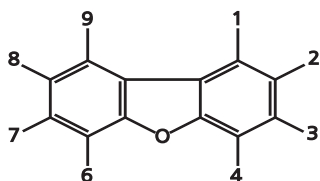
NATIVE CHLORINATED DIBENZOFURANS (PCDFs)

| Catalogue Number | Product (toluene or nonane solution) | Qty/Conc |
|------------------|--|-----------------|
| DF-1348-S | 1,3,4,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1349-S | 1,3,4,9-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1367-S | 1,3,6,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1368-S | 1,3,6,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1369-S | 1,3,6,9-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1378-S | 1,3,7,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1467-S | 1,4,6,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1478-S | 1,4,7,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2346-S | 2,3,4,6-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2347-S | 2,3,4,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2348-S | 2,3,4,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2368-S | 2,3,6,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2378-S | 2,3,7,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2467-S | 2,4,6,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-2468-S | 2,4,6,8-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-3467-S | 3,4,6,7-Tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12347-S | 1,2,3,4,7-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12348-S | 1,2,3,4,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12367-S | 1,2,3,6,7-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12378-S | 1,2,3,7,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12379-S | 1,2,3,7,9-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12389-S | 1,2,3,8,9-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12467-S | 1,2,4,6,7-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12468-S | 1,2,4,6,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12478-S | 1,2,4,7,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-13467-S | 1,3,4,6,7-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-13478-S | 1,3,4,7,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-13479-S | 1,3,4,7,9-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-13678-S | 1,3,6,7,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-23467-S | 2,3,4,6,7-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-23469-S | 2,3,4,6,9-Pentachlorodibenzofuran (or 1,4,6,7,8) | 1.2 ml 50 µg/ml |
| DF-23478-S | 2,3,4,7,8-Pentachlorodibenzofuran | 1.2 ml 50 µg/ml |



NATIVE CHLORINATED DIBENZOFURANS (PCDFs)

| Catalogue Number | Product (toluene or nonane solution) | Qty/Conc |
|------------------|---------------------------------------|-----------------|
| DF-123467-S | 1,2,3,4,6,7-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-123468-S | 1,2,3,4,6,8-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-123478-S | 1,2,3,4,7,8-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-123489-S | 1,2,3,4,8,9-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-123678-S | 1,2,3,6,7,8-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-123789-S | 1,2,3,7,8,9-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-124678-S | 1,2,4,6,7,8-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-234678-S | 2,3,4,6,7,8-Hexachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1234678-S | 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1234689-S | 1,2,3,4,6,8,9-Heptachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-1234789-S | 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 1.2 ml 50 µg/ml |
| DF-12346789-S | Octachlorodibenzofuran | 1.2 ml 50 µg/ml |

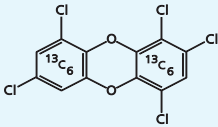
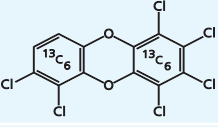
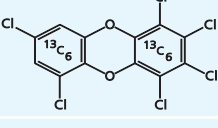
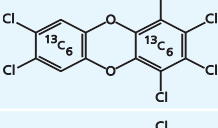
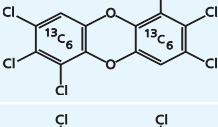
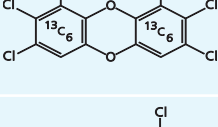
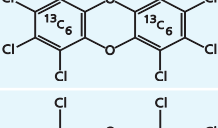
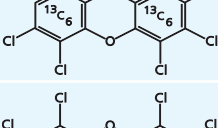
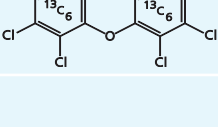
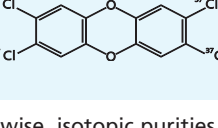


MASS-LABELLED CHLORINATED DIBENZO-p-DIOXINS

| Catalogue Number | Product |
|------------------|---|
| MDD-0 |  <p>[¹³C₁₂]Dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-2 |  <p>2-Chloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-23 |  <p>2,3-Dichloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-237 |  <p>2,3,7-Trichloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1234 |  <p>1,2,3,4-Tetrachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1278 |  <p>1,2,7,8-Tetrachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1368 |  <p>1,3,6,8-Tetrachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1378 |  <p>1,3,7,8-Tetrachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-2378 |  <p>2,3,7,8-Tetrachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-12378 |  <p>1,2,3,7,8-Pentachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-12389 |  <p>1,2,3,8,9-Pentachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-12478 |  <p>1,2,4,7,8-Pentachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

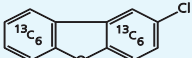
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED DIBENZO-p-DIOXINS

| Catalogue Number | Product |
|------------------|---|
| MDD-12479 |  <p>1,2,4,7,9-Pentachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-123467 |  <p>1,2,3,4,6,7-Hexachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-123468 |  <p>1,2,3,4,6,8-Hexachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-123478 |  <p>1,2,3,4,7,8-Hexachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-123678 |  <p>1,2,3,6,7,8-Hexachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-123789 |  <p>1,2,3,7,8,9-Hexachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1234678 |  <p>1,2,3,4,6,7,8-Heptachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-1234679 |  <p>1,2,3,4,6,7,9-Heptachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDD-12346789 |  <p>Octachloro[¹³C₁₂]dibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MCDD-2378 |  <p>2,3,7,8-[³⁷Cl₄]-Tetrachlorodibenzo-p-dioxin 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene (isotopic purity; 94 to 95%)</p> |

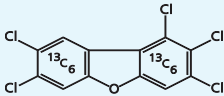
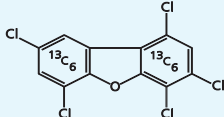
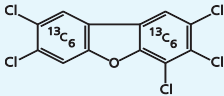
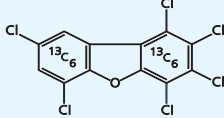
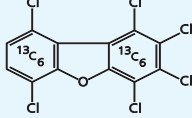
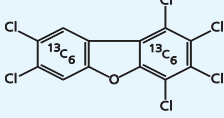
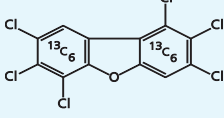
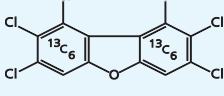
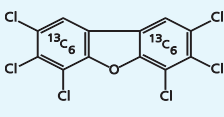
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED DIBENZOFURANS

| Catalogue Number | Product |
|------------------|---|
| MDF-0 |  <p>[¹³C₁₂]Dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-2 |  <p>2-Chloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-23 |  <p>2,3-Dichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-238 |  <p>2,3,8-Trichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-1234 |  <p>1,2,3,4-Tetrachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-1278 |  <p>1,2,7,8-Tetrachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-1368 |  <p>1,3,6,8-Tetrachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-2378 |  <p>2,3,7,8-Tetrachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-12346 |  <p>1,2,3,4,6-Pentachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

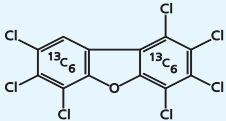
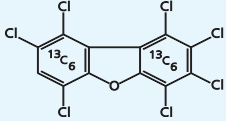
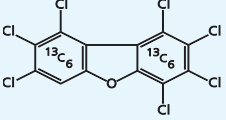
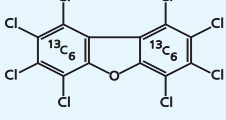
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED DIBENZOFURANS

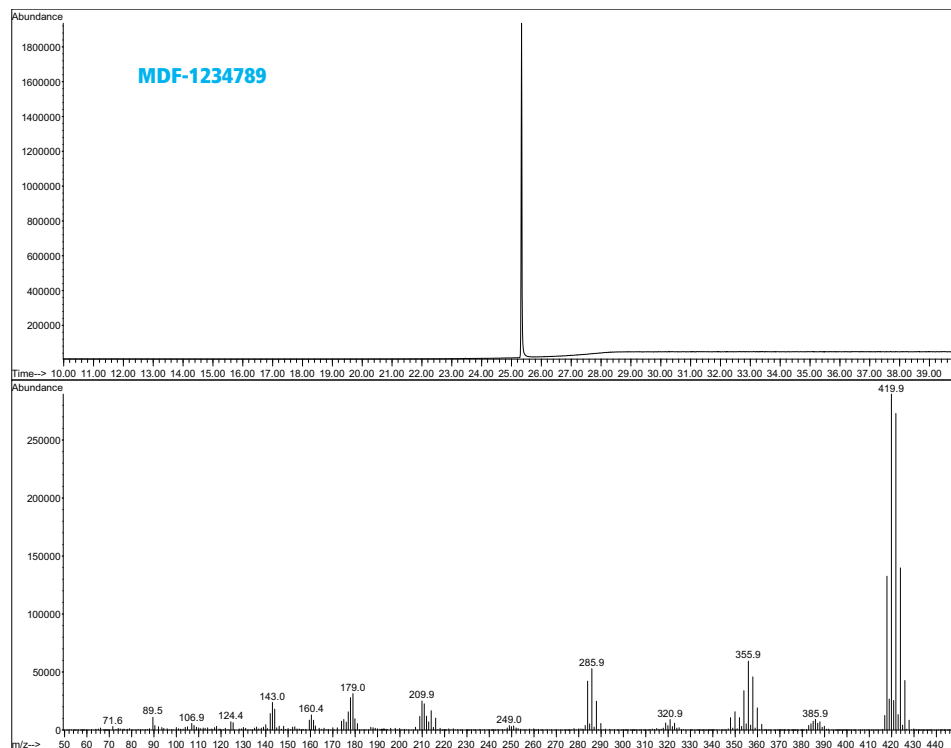
| Catalogue Number | Product |
|-------------------|--|
| MDF-12378 |  <p>1,2,3,7,8-Pentachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-13468 |  <p>1,3,4,6,8-Pentachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-23478 |  <p>2,3,4,7,8-Pentachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-123468 |  <p>1,2,3,4,6,8-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-123469 |  <p>1,2,3,4,6,9-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-123478 |  <p>1,2,3,4,7,8-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-123678 |  <p>1,2,3,6,7,8-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-123789 |  <p>1,2,3,7,8,9-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-234678 |  <p>2,3,4,6,7,8-Hexachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

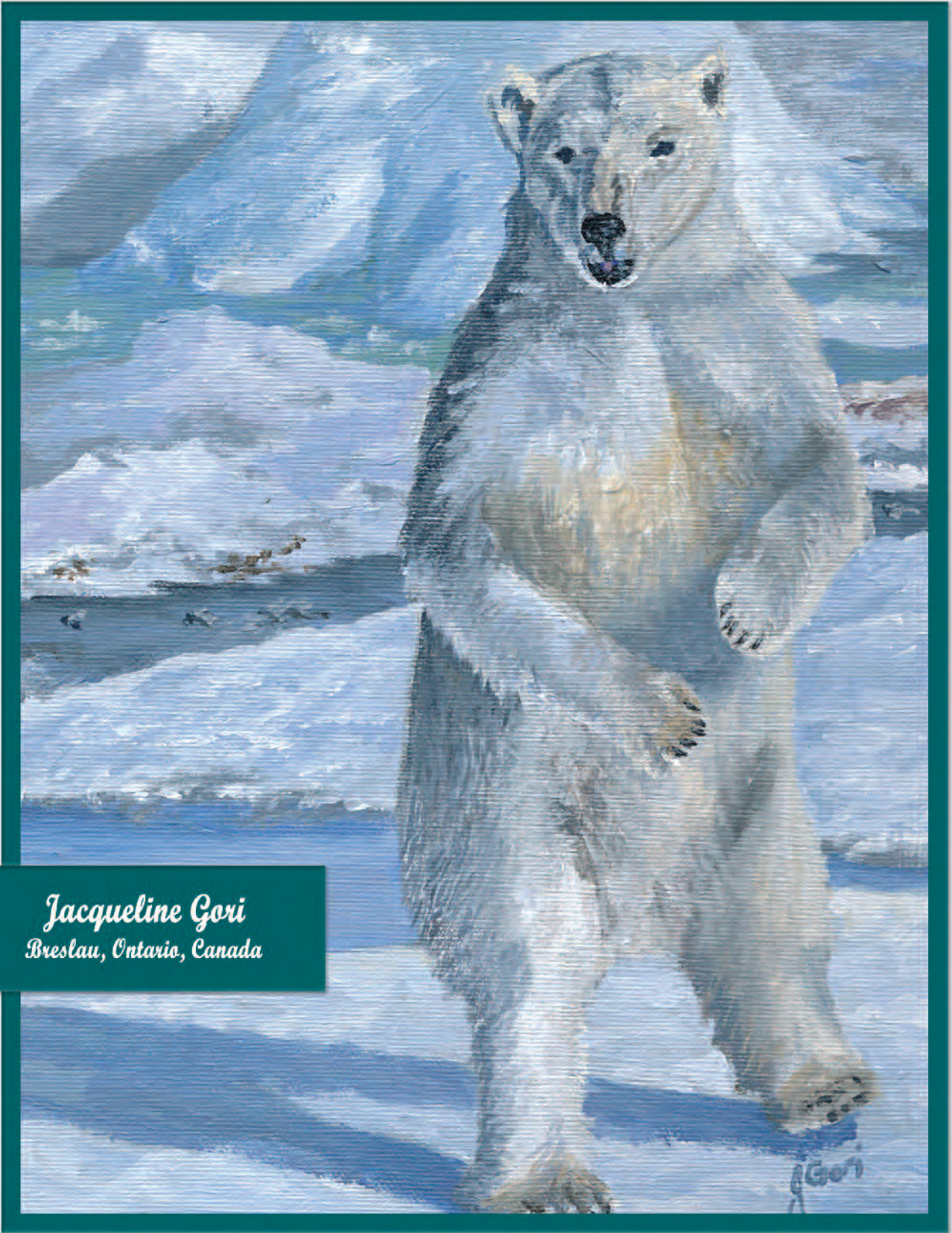
MASS-LABELLED CHLORINATED DIBENZOFURANS

| Catalogue Number | Product |
|------------------|---|
| MDF-1234678 |  <p>1,2,3,4,6,7,8-Heptachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-1234689 |  <p>1,2,3,4,6,8,9-Heptachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-1234789 |  <p>1,2,3,4,7,8,9-Heptachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MDF-12346789 |  <p>Octachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.



HRGC/LRMS data of 1,2,3,4,7,8,9-Heptachloro[¹³C₁₂]dibenzofuran (30m DB-5 column).



Jacqueline Gori
Breslau, Ontario, Canada

PCDDs & PCDFs:

JAPANESE INDUSTRIAL STANDARD METHODS JIS K 0311 AND JIS K 0312 & ADDITIONAL PCDD/PCDF SPECIALTY SOLUTION/MIXTURES

Wellington has also prepared several other calibration sets containing native and mass-labelled chlorinated dibenzo-p-dioxins (PCDDs) and chlorinated dibenzofurans (PCDFs). Some of these have become quite popular as they are multi-point sets including as many as 13 individual calibration solutions. This allows some flexibility in selecting the calibration range. In addition, these calibration sets and their support solutions incorporate additional mass-labelled PCDDs and PCDFs which can be used as part of a more involved sampling and sample processing regime.

All of these calibration sets and support solutions are suitable for use with the Japanese Industrial Standards JIS K 0311 and JIS K 0312. In addition, provided that all performance criteria are achieved, they can be used with other regulatory methods from other countries.

In our previous catalogue, we introduced DFP-CVS-B10 which is a combination calibration set for PCDD, PCDF and dioxin-like PCB congener analysis. These solutions remain very popular and will continue to be offered.



DF-CVS-A10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------------|------------------------------------|-----------------------------|
| DF-CVS-A10-Set 1 | CS1/CS3/CS5/CS7/CS9 | 1 kit (5 x 200 µl ampoules) |
| DF-CVS-A10-Set 2 | CS2/CS4/CS6/CS8/CS10 | 1 kit (5 x 200 µl ampoules) |
| DF-CVS-A10-Set 3 | CS3/CS5/CS7/CS9/CS11 | 1 kit (5 x 200 µl ampoules) |
| DF-A10-CSL | CSL Extended Calibration/Low Level | 200 µl |
| DF-A10-CS1 | CS1 | 200 µl |
| DF-A10-CS2 | CS2 | 200 µl |
| DF-A10-CS3 | CS3 | 200 µl |
| DF-A10-CS4 | CS4 | 200 µl |

| | DF-A10-CSL (ng/ml) | DF-A10-CS1 (ng/ml) | DF-A10-CS2 (ng/ml) | DF-A10-CS3 (ng/ml) | DF-A10-CS4 (ng/ml) |
|--|-----------------------|-----------------------|-----------------------|-----------------------|-----------------------|
| NATIVE PCDDs & PCDFs | | | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| Octachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.4 | 1 | 2 |
| | | | | | |
| 1,3,6,8-Tetrachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,7,8-Tetrachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,8,9-Tetrachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| Octachlorodibenzofuran | 0.1 | 0.2 | 0.4 | 1 | 2 |
| | | | | | |
| MASS-LABELLED PCDDs & PCDFs | | | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 |
| | | | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|-------------------------------------|----------|
| DF-A10-CS5 | CS5 | 200 µl |
| DF-A10-CS6 | CS6 | 200 µl |
| DF-A10-CS7 | CS7 | 200 µl |
| DF-A10-CS8 | CS8 | 200 µl |
| DF-A10-CS9 | CS9 | 200 µl |
| DF-A10-CS10 | CS10 | 200 µl |
| DF-A10-CS11 | CS11 | 200 µl |
| DF-A10-CSH | CSH Extended Calibration/High Level | 200 µl |

| DF-A10-CS5 (ng/ml) | DF-A10-CS6 (ng/ml) | DF-A10-CS7 (ng/ml) | DF-A10-CS8 (ng/ml) | DF-A10-CS9 (ng/ml) | DF-A10-CS10 (ng/ml) | DF-A10-CS11 (ng/ml) | DF-A10-CSH (ng/ml) |
|--------------------|--------------------|--------------------|--------------------|--------------------|---------------------|---------------------|--------------------|
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 | 2000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 | 1000 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 | 2000 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |

DF-CVS-B10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------------|---------------------------|-----------------------------|
| DF-CVS-B10-Set 1 | CS1/CS3/CS5/CS7/CS9 | 1 kit (5 x 200 µl ampoules) |
| DF-CVS-B10-Set 2 | CS2/CS4/CS6/CS8/CS10 | 1 kit (5 x 200 µl ampoules) |
| DF-CVS-B10-Set 3 | CS3/CS5/CS7/CS9/CS11 | 1 kit (5 x 200 µl ampoules) |
| DF-B10-CS1 | CS1 | 200 µl |
| DF-B10-CS2 | CS2 | 200 µl |
| DF-B10-CS3 | CS3 | 200 µl |
| DF-B10-CS4 | CS4 | 200 µl |

| | DF-B10-CS1 (ng/ml) | DF-B10-CS2 (ng/ml) | DF-B10-CS3 (ng/ml) | DF-B10-CS4 (ng/ml) |
|--|-----------------------|-----------------------|-----------------------|-----------------------|
| NATIVE PCDDs & PCDFs | | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.5 | 1 |
| 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.2 | 0.4 | 1 | 2 |
| Octachlorodibenzo-p-dioxin | 0.5 | 1 | 2.5 | 5 |
| 1,3,6,8-Tetrachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,7,8-Tetrachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,8,9-Tetrachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.1 | 0.2 | 0.5 | 1 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.2 | 0.4 | 1 | 2 |
| Octachlorodibenzofuran | 0.5 | 1 | 2.5 | 5 |
| MASS-LABELLED PCDDs & PCDFs | | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------|----------|
| DF-B10-CS5 | CS5 | 200 µl |
| DF-B10-CS6 | CS6 | 200 µl |
| DF-B10-CS7 | CS7 | 200 µl |
| DF-B10-CS8 | CS8 | 200 µl |
| DF-B10-CS9 | CS9 | 200 µl |
| DF-B10-CS10 | CS10 | 200 µl |
| DF-B10-CS11 | CS11 | 200 µl |

| DF-B10-CS5 (ng/ml) | DF-B10-CS6 (ng/ml) | DF-B10-CS7 (ng/ml) | DF-B10-CS8 (ng/ml) | DF-B10-CS9 (ng/ml) | DF-B10-CS10 (ng/ml) | DF-B10-CS11 (ng/ml) |
|--------------------|--------------------|--------------------|--------------------|--------------------|---------------------|---------------------|
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 10 | 25 | 50 | 100 | 250 | 500 | 1000 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 2 | 5 | 10 | 20 | 50 | 100 | 200 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 4 | 10 | 20 | 40 | 100 | 200 | 400 |
| 10 | 25 | 50 | 100 | 250 | 500 | 1000 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 20 | 20 | 20 | 20 | 20 | 20 | 20 |

DF-CVS-C10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------|--|-----------------------------|
| DF-CVS-C10 | PCDD & PCDF Calibration Solutions CS1 / CS2 / CS3 / CS4 / CS5 / CS6 / CS7 | 1 kit (7 x 200 µl ampoules) |
| DF-C10-CS1 | CS1 | 200 µl |
| DF-C10-CS2 | CS2 | 200 µl |
| DF-C10-CS3 | CS3 | 200 µl |

| | DF-C10-CS1 (ng/ml) | DF-C10-CS2 (ng/ml) | DF-C10-CS3 (ng/ml) |
|--|-----------------------|-----------------------|-----------------------|
| NATIVE PCDDs & PCDFs | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| Octachlorodibenzo-p-dioxin | 0.5 | 2.5 | 10 |
| | | | |
| 1,3,6,8-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,8,9-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.2 | 1 | 4 |
| Octachlorodibenzofuran | 0.5 | 2.5 | 10 |
| EXTRACTION SPIKE: LABELLED PCDDs & PCDFs | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 |
| | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 |
| SYRINGE SPIKE: LABELLED PCDDs | | | |
| 1,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,4,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,7,9-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |
| SAMPLING SPIKE | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------|----------|
| DF-C10-CS4 | CS4 | 200 µl |
| DF-C10-CS5 | CS5 | 200 µl |
| DF-C10-CS6 | CS6 | 200 µl |
| DF-C10-CS7 | CS7 | 200 µl |

*FOR SUPPORT SOLUTIONS, see DF-ST-C, DF-LCS-B, DF-IS-J, and DF-SS-A (and their dilutions).

| | DF-C10-CS4 (ng/ml) | DF-C10-CS5 (ng/ml) | DF-C10-CS6 (ng/ml) | DF-C10-CS7 (ng/ml) |
|--|-----------------------|-----------------------|-----------------------|-----------------------|
| NATIVE PCDDs & PCDFs | | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| Octachlorodibenzo-p-dioxin | 50 | 250 | 1000 | 2500 |
| | | | | |
| 1,3,6,8-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,8,9-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 2,3,7,8-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| Octachlorodibenzofuran | 50 | 250 | 1000 | 2500 |
| EXTRACTION SPIKE: LABELLED PCDDs & PCDFs | | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 |
| | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 |
| SYRINGE SPIKE: LABELLED PCDDs | | | | |
| 1,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,4,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,9-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| SAMPLING SPIKE | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 | 10 | 10 | 10 |

DFP-CVS-B10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|-----------------------------|
| DFP-CVS-B10 | PCDD/PCDF/PCB Calibration Solutions CS1 / CS2 / CS3 / CS4 / CS5 / CS6 / CS7 | 1 kit (7 x 200 µl ampoules) |
| DFP-B10-CS1 | CS1 | 200 µl |
| DFP-B10-CS2 | CS2 | 200 µl |
| DFP-B10-CS3 | CS3 | 200 µl |

| | DFP-B10-CS1 (ng/ml) | DFP-B10-CS2 (ng/ml) | DFP-B10-CS3 (ng/ml) |
|--|------------------------|------------------------|------------------------|
| NATIVE PCDDs & PCDFs | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.1 | 0.5 | 2 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.2 | 1 | 4 |
| Octachlorodibenzo-p-dioxin | 0.5 | 2.5 | 10 |
| 1,3,6,8-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,8,9-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.1 | 0.5 | 2 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.2 | 1 | 4 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.2 | 1 | 4 |
| Octachlorodibenzofuran | 0.5 | 2.5 | 10 |
| NATIVE PCBs | | | |
| | IUPAC | | |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.2 | 1 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.2 | 1 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.2 | 1 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.2 | 1 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.2 | 1 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.2 | 1 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.2 | 1 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.2 | 1 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.2 | 1 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.2 | 1 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.2 | 1 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 0.2 | 1 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 0.2 | 1 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.2 | 1 |
| EXTRACTION SPIKE: 13C PCDDs/PCDFs/PCBs | | | |
| 1,3,6,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| Octachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 20 | 20 | 20 |
| 2,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 |
| Octachloro[¹³ C] ₁₂ dibenzofuran | 20 | 20 | 20 |
| 3,3',4,4'-Tetrachloro[¹³ C] ₁₂ biphenyl | 77L | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C] ₁₂ biphenyl | 81L | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C] ₁₂ biphenyl | 105L | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 114L | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 118L | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 123L | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 126L | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C] ₁₂ biphenyl | 156L | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C] ₁₂ biphenyl | 157L | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C] ₁₂ biphenyl | 167L | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C] ₁₂ biphenyl | 169L | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C] ₁₂ biphenyl | 170L | 10 | 10 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C] ₁₂ biphenyl | 180L | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C] ₁₂ biphenyl | 189L | 10 | 10 |
| SYRINGE SPIKE: 13C PCDDs & PCBs | | | |
| 1,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,4,7,8-Pentachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 1,2,3,4,6,7,9-Heptachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 2,3',4',5-Tetrachloro[¹³ C] ₁₂ biphenyl | 70L | 10 | 10 |
| 2,3,3',5,5'-Pentachloro[¹³ C] ₁₂ biphenyl | 111L | 10 | 10 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C] ₁₂ biphenyl | 138L | 10 | 10 |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C] ₁₂ biphenyl | 178L | 10 | 10 |
| SAMPLING SPIKE | | | |
| 1,2,3,4-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 |
| 3,3',4,5'-Tetrachloro[¹³ C] ₁₂ biphenyl | 79L | 10 | 10 |

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|---------------------------|------------------------|------------------------|------------------------|
| DFP-B10-CS4 | CS4 | 200 µl | | |
| DFP-B10-CS5 | CS5 | 200 µl | | |
| DFP-B10-CS6 | CS6 | 200 µl | | |
| DFP-B10-CS7 | CS7 | 200 µl | | |
| *FOR PCDD/PCDF/PCB SUPPORT SOLUTIONS, see DFP-LCS-B (page 59), DFP-IS-B10 (page 61), and DFP-SS-A10 (page 61). | | | | |
| *FOR PCDD/PCDF SUPPORT SOLUTIONS, see DF-ST-C, DF-LCS-B, DF-IS-J, and DF-SS-A. | | | | |
| *FOR NATIVE AND MASS-LABELLED PCB SUPPORT SOLUTIONS, see pages 92 and 94. | | | | |
| | DFP-B10-CS4 (ng/ml) | DFP-B10-CS5 (ng/ml) | DFP-B10-CS6 (ng/ml) | DFP-B10-CS7 (ng/ml) |
| NATIVE PCDDs & PCDFs | | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 10 | 50 | 200 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 20 | 100 | 400 | 1000 |
| Octachlorodibenzo-p-dioxin | 50 | 250 | 1000 | 2500 |
| 1,3,6,8-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,8,9-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 2,3,7,8-Tetrachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 10 | 50 | 200 | 500 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 20 | 100 | 400 | 1000 |
| Octachlorodibenzofuran | 50 | 250 | 1000 | 2500 |
| NATIVE PCBs | | | | |
| | IUPAC | | | |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 20 | 400 | 1000 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 20 | 400 | 1000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 20 | 400 | 1000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 20 | 400 | 1000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 20 | 400 | 1000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 20 | 400 | 1000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 20 | 400 | 1000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 20 | 400 | 1000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 20 | 400 | 1000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 20 | 400 | 1000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 20 | 400 | 1000 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 20 | 400 | 1000 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 20 | 400 | 1000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 20 | 400 | 1000 |
| EXTRACTION SPIKE: 13C PCDDs/PCDFs/PCBs | | | | |
| 1,3,6,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 2,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 20 | 20 | 20 | 20 |
| 2,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-Pentachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-Pentachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-Hexachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C] ₁₂ dibenzofuran | 10 | 10 | 10 | 10 |
| Octachloro[¹³ C] ₁₂ dibenzofuran | 20 | 20 | 20 | 20 |
| 3,3',4,4'-Tetrachloro[¹³ C] ₁₂ biphenyl | 77L | 10 | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C] ₁₂ biphenyl | 81L | 10 | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C] ₁₂ biphenyl | 105L | 10 | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 114L | 10 | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 118L | 10 | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 123L | 10 | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C] ₁₂ biphenyl | 126L | 10 | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C] ₁₂ biphenyl | 156L | 10 | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C] ₁₂ biphenyl | 157L | 10 | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C] ₁₂ biphenyl | 167L | 10 | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C] ₁₂ biphenyl | 169L | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C] ₁₂ biphenyl | 170L | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C] ₁₂ biphenyl | 180L | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C] ₁₂ biphenyl | 189L | 10 | 10 | 10 |
| SYRINGE SPIKE: 13C PCDDs & PCBs | | | | |
| 1,3,7,8-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,4,7,8-Pentachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8-Hexachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,9-Heptachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 2,3',4',5-Tetrachloro[¹³ C] ₁₂ biphenyl | 70L | 10 | 10 | 10 |
| 2,3,3',5,5'-Pentachloro[¹³ C] ₁₂ biphenyl | 111L | 10 | 10 | 10 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C] ₁₂ biphenyl | 138L | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C] ₁₂ biphenyl | 178L | 10 | 10 | 10 |
| SAMPLING SPIKE | | | | |
| 1,2,3,4-Tetrachloro[¹³ C] ₁₂ dibenzo-p-dioxin | 10 | 10 | 10 | 10 |
| 3,3',4,5'-Tetrachloro[¹³ C] ₁₂ biphenyl | 79L | 10 | 10 | 10 |

NK-CVS-J

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------------|
| NK-CVS-J | PCDD & PCDF Calibration Solutions CS1-J, CS2-J, CS3-J, CS4-J, CS5-J and CS6-J | 1 kit (6 ampoules) |
| NK-CS1-J | CS1-J | 200 µl |
| NK-CS2-J | CS2-J | 200 µl |
| NK-CS3-J | CS3-J | 200 µl |
| NK-CS4-J | CS4-J | 200 µl |
| NK-CS5-J | CS5-J | 200 µl |
| NK-CS6-J | CS6-J | 200 µl |

| | NK-CS1-J (ng/ml) | NK-CS2-J (ng/ml) | NK-CS3-J (ng/ml) | NK-CS4-J (ng/ml) | NK-CS5-J (ng/ml) | NK-CS6-J (ng/ml) |
|--|---------------------|---------------------|---------------------|---------------------|---------------------|---------------------|
| NATIVE PCDDs & PCDFs | | | | | | |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| Octachlorodibenzo-p-dioxin | 0.5 | 2.0 | 10 | 50 | 250 | 500 |
| | | | | | | |
| 1,2,7,8-Tetrachlorodibenzofuran | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 2,3,7,8-Tetrachlorodibenzofuran | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 0.1 | 0.4 | 2 | 10 | 50 | 100 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 0.2 | 0.8 | 4 | 20 | 100 | 200 |
| Octachlorodibenzofuran | 0.5 | 2.0 | 10 | 50 | 250 | 500 |
| | | | | | | |
| MASS-LABELLED PCDDs & PCDFs* | | | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 40 | 40 | 40 | 40 | 40 | 40 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 40 | 40 | 40 | 40 | 40 | 40 |
| | | | | | | |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 20 | 20 | 20 | 20 | 20 | 20 |
| | | | | | | |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 20 | 20 | 20 | 20 | 20 | 20 |

* Support Solutions for **NK-CVS-J** see: **NK-LCS-T**, **NK-IS-J4** and **NK-IS-J5**.

Support solutions for **DF-CVS-A10** and **DF-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|----------|
| DF-LCS-A | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-A200 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-A40 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |

| | DF-LCS-A (ng/ml) | DF-LCS-A200 (ng/ml) | DF-LCS-A40 (ng/ml) |
|--|---------------------|------------------------|-----------------------|
| MASS-LABELLED PCDDs | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 2000 | 400 | 80 |
| MASS-LABELLED PCDFs | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 2000 | 400 | 80 |

DF-LCS-B

Support solutions for **DF-CVS-A10**, **DF-CVS-B10**, **DF-CVS-C10**, and **DFP-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|----------|
| DF-LCS-B | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-B200 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-B40 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |

| | DF-LCS-B (ng/ml) | DF-LCS-B200 (ng/ml) | DF-LCS-B40 (ng/ml) |
|--|---------------------|------------------------|-----------------------|
| MASS-LABELLED PCDDs | | | |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 2000 | 400 | 80 |
| MASS-LABELLED PCDFs | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 2000 | 400 | 80 |

Support solutions for **DF-CVS-A10**, **DF-CVS-B10**, **DF-CVS-C10**, and **DFP-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|----------|
| DF-LCS-C | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-C200 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-LCS-C40 | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |

| | DF-LCS-C (ng/ml) | DF-LCS-C200 (ng/ml) | DF-LCS-C40 (ng/ml) |
|--|---------------------|------------------------|-----------------------|
| MASS-LABELLED PCDDs | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 2000 | 400 | 80 |
| MASS-LABELLED PCDFs | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | 1000 | 200 | 40 |
| Octachloro[¹³ C ₁₂]dibenzofuran | 2000 | 400 | 80 |

MASS-LABELLED PCDDs & PCDFs: SOLUTION/MIXTURES

NOTE CONCENTRATIONS: DS-1000 and FS-1000 = µg/ml
NK-LCS-O and NK-LCS-T = ng/ml

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|----------|
| DS-1000 | Mass-Labelled PCDD Solution/Mixture | 1.2 ml |
| FS-1000 | Mass-Labelled PCDF Solution/Mixture | 1.2 ml |
| NK-LCS-O | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |
| NK-LCS-T | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |

| | DS-1000 (µg/ml) | FS-1000 (µg/ml) | NK-LCS-O (ng/ml) | NK-LCS-T (ng/ml) |
|--|--------------------|--------------------|---------------------|---------------------|
| MASS-LABELLED PCDDs | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | — | 100 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | — | 100 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | — | 100 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | — | 100 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | — | — | — |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 1.0 | — | 100 | 40 |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | 2.0 | — | 200 | 80 |
| MASS-LABELLED PCDFs | | | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | — |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | — |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | — | 1.0 | 100 | 40 |
| Octachloro[¹³ C ₁₂]dibenzofuran | — | — | 200 | 80 |

MASS-LABELLED PCDDs & PCDFs: SOLUTION/MIXTURES

Support solutions for **DF-CVS-A10** and **DF-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|--|----------------|-------------------|------------------|
| DF-IS-A | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| DF-IS-A200 | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| DF-IS-A40 | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-A | DF-IS-A200 | DF-IS-A40 |
| MASS-LABELLED PCDD | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | | 1.0 | 200 | 40 |
| DF-IS-B | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-B200 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-B40 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-B | DF-IS-B200 | DF-IS-B40 |
| MASS-LABELLED PCDD/PCDF | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | | 1.0 | 200 | 40 |
| 1,2,3,4-Tetrachloro ^[13C₁₂] dibenzofuran | | 1.0 | 200 | 40 |
| DF-IS-C | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-C200 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-C40 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-C | DF-IS-C200 | DF-IS-C40 |
| MASS-LABELLED PCDD/PCDF | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | | 1.0 | 200 | 40 |
| 1,2,7,8-Tetrachloro ^[13C₁₂] dibenzofuran | | 1.0 | 200 | 40 |

MASS-LABELLED PCDDs & PCDFs: SOLUTION/MIXTURES

Support solutions for **DF-CVS-A10** and **DF-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|--|----------------|-------------------|------------------|
| DF-IS-D | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-D200 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-D40 | Mass-Labelled PCDD/PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-D | DF-IS-D200 | DF-IS-D40 |
| MASS-LABELLED PCDD/PCDF | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1.0 | 200 | 40 |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| DF-IS-E | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-E200 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-E40 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-E | DF-IS-E200 | DF-IS-E40 |
| MASS-LABELLED PCDFs | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,3,6,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| DF-IS-F | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-F200 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-F40 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-F | DF-IS-F200 | DF-IS-F40 |
| MASS-LABELLED PCDFs | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |

MASS-LABELLED PCDDs & PCDFs: SOLUTION/MIXTURES

Support solutions for **DF-CVS-A10** and **DF-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|---|----------------|-------------------|------------------|
| DF-IS-G | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-G200 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-G40 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-G | DF-IS-G200 | DF-IS-G40 |
| MASS-LABELLED PCDFs | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4,6-Pentachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| DF-IS-H | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-H200 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-H40 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-H | DF-IS-H200 | DF-IS-H40 |
| MASS-LABELLED PCDFs | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| DF-IS-I | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-I200 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| DF-IS-I40 | Mass-Labelled PCDF Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-I | DF-IS-I200 | DF-IS-I40 |
| MASS-LABELLED PCDFs | | (µg/ml) | (ng/ml) | (ng/ml) |
| 1,2,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6-Pentachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1.0 | 200 | 40 |

MASS-LABELLED PCDDs & PCDFs: SOLUTION/MIXTURES

Support solutions for **DFP-CVS-B10** and **DF-CVS-C10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|---|----------------|-------------------|------------------|
| DF-IS-J | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| DF-IS-J100 | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| DF-IS-J20 | Mass-Labelled PCDD Internal Standard Solution | 1.2 ml | | |
| | | DF-IS-J | DF-IS-J100 | DF-IS-J20 |
| MASS-LABELLED PCDDs | | (ng/ml) | (ng/ml) | (ng/ml) |
| 1,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,4,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,4,6,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,4,6,7,9-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1000 | 100 | 20 |
| | | DF-IS-J | DF-IS-J100 | DF-IS-J20 |
| DF-SS-A | Mass-Labelled PCDD Sampling Spike Solution | 1.2 ml | | |
| DF-SS-A100 | Mass-Labelled PCDD Sampling Spike Solution | 1.2 ml | | |
| DF-SS-A20 | Mass-Labelled PCDD Sampling Spike Solution | 1.2 ml | | |
| | | DF-SS-A | DF-SS-A100 | DF-SS-A20 |
| MASS-LABELLED PCDD | | (ng/ml) | (ng/ml) | (ng/ml) |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1000 | 100 | 20 |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|----------------------------|----------|
| MDF-1278-1 | | 1.2 ml |
| 1,2,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 1 µg/ml |
| MDF-12346-1 | | 1.2 ml |
| 1,2,3,4,6-Pentachloro[¹³ C ₁₂]dibenzofuran | | 1 µg/ml |
| MDF-123469-1 | | 1.2 ml |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | | 1 µg/ml |
| MDF-1234689-1 | | 1.2 ml |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 1 µg/ml |
| NK-IS-J4 | Internal Standard Solution | 1.2 ml |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 40 ng/ml |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 40 ng/ml |
| NK-IS-J5 | Internal Standard Solution | 1.2 ml |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | | 40 ng/ml |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | | 40 ng/ml |

MASS-LABELLED PCDDs/PCDFs/PCBs: SOLUTION/MIXTURES

Support solutions for **DFP-CVS-B10**, **DF-CVS-C10**, and **PCB-CVS-B10** (see page 90)

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--|---|-----------|--------------|-------------|
| DFP-LCS-B | Mass-Labelled PCDDs/PCDFs/PCBs Solution/Mixture | 1.2 ml | | |
| DFP-LCS-B100 | Mass-Labelled PCDDs/PCDFs/PCBs Solution/Mixture | 1.2 ml | | |
| DFP-LCS-B20 | Mass-Labelled PCDDs/PCDFs/PCBs Solution/Mixture | 1.2 ml | | |
| | | DFP-LCS-B | DFP-LCS-B100 | DFP-LCS-B20 |
| MASS-LABELLED PCDDs & PCDFs | | (ng/ml) | (ng/ml) | (ng/ml) |
| 1,3,6,8-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 2,3,7,8-Tetrachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,7,8-Pentachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,4,7,8-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,6,7,8-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,7,8,9-Hexachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| 1,2,3,4,6,7,8-Heptachloro ^[13C₁₂] dibenzo-p-dioxin | | 1000 | 100 | 20 |
| Octachloro ^[13C₁₂] dibenzo-p-dioxin | | 2000 | 200 | 40 |
| 2,3,7,8-Tetrachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,7,8-Pentachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 2,3,4,7,8-Pentachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,4,7,8-Hexachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,6,7,8-Hexachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,7,8,9-Hexachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 2,3,4,6,7,8-Hexachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,4,6,7,8-Heptachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| 1,2,3,4,7,8,9-Heptachloro ^[13C₁₂] dibenzofuran | | 1000 | 100 | 20 |
| Octachloro ^[13C₁₂] dibenzofuran | | 2000 | 200 | 40 |
| MASS-LABELLED PCBs | IUPAC | | | |
| 3,3',4,4'-Tetrachloro ^[13C₁₂] biphenyl | 77L | 1000 | 100 | 20 |
| 3,4,4',5-Tetrachloro ^[13C₁₂] biphenyl | 81L | 1000 | 100 | 20 |
| 2,3,3',4,4'-Pentachloro ^[13C₁₂] biphenyl | 105L | 1000 | 100 | 20 |
| 2,3,4,4',5-Pentachloro ^[13C₁₂] biphenyl | 114L | 1000 | 100 | 20 |
| 2,3',4,4',5-Pentachloro ^[13C₁₂] biphenyl | 118L | 1000 | 100 | 20 |
| 2',3,4,4',5-Pentachloro ^[13C₁₂] biphenyl | 123L | 1000 | 100 | 20 |
| 3,3',4,4',5-Pentachloro ^[13C₁₂] biphenyl | 126L | 1000 | 100 | 20 |
| 2,3,3',4,4',5-Hexachloro ^[13C₁₂] biphenyl | 156L | 1000 | 100 | 20 |
| 2,3,3',4,4',5'-Hexachloro ^[13C₁₂] biphenyl | 157L | 1000 | 100 | 20 |
| 2,3',4,4',5,5'-Hexachloro ^[13C₁₂] biphenyl | 167L | 1000 | 100 | 20 |
| 3,3',4,4',5,5'-Hexachloro ^[13C₁₂] biphenyl | 169L | 1000 | 100 | 20 |
| 2,2',3,3',4,4',5-Heptachloro ^[13C₁₂] biphenyl | 170L | 1000 | 100 | 20 |
| 2,2',3,4,4',5,5'-Heptachloro ^[13C₁₂] biphenyl | 180L | 1000 | 100 | 20 |
| 2,3,3',4,4',5,5'-Heptachloro ^[13C₁₂] biphenyl | 189L | 1000 | 100 | 20 |

MASS-LABELLED PCDDs/PCDFs/PCBs: SOLUTION/MIXTURES

These three solutions were designed and prepared as support solutions to be used with the following calibration sets:

DF-CVS-A10

DF-CVS-B10

as well as:

PCB-CVS-A10 (see Page 88)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--|--|----------|
| DFP-LCS-A | Mass-Labelled PCDD/PCDF/PCB Solution/Mixture | 1.2 ml |
| MASS-LABELLED PCDDs | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 10 ng/ml |
| Octachloro[¹³ C ₁₂]dibenzo-p-dioxin | | 20 ng/ml |
| MASS-LABELLED PCDFs | | |
| 2,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 2,3,4,7,8-Pentachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,4,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,7,8,9-Hexachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 2,3,4,6,7,8-Hexachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,4,6,7,8-Heptachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| 1,2,3,4,7,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | | 10 ng/ml |
| Octachloro[¹³ C ₁₂]dibenzofuran | | 20 ng/ml |
| MASS-LABELLED PCBs | | |
| | IUPAC | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 ng/ml |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 ng/ml |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 10 ng/ml |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 10 ng/ml |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 10 ng/ml |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 10 ng/ml |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 ng/ml |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 10 ng/ml |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 10 ng/ml |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 10 ng/ml |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 ng/ml |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 ng/ml |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 10 ng/ml |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 10 ng/ml |
| DFP-IS-A | Mass-Labelled PCDF/PCB Syringe Spike | 1.2 ml |
| | IUPAC | |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 ng/ml |
| 1,2,3,4,6,9-Hexachloro[¹³ C ₁₂]dibenzofuran | — | 10 ng/ml |
| 1,2,3,4,6,8,9-Heptachloro[¹³ C ₁₂]dibenzofuran | — | 10 ng/ml |
| DFP-SS-A | Mass-Labelled PCDD/PCB Sampling Spike | 1.2 ml |
| | IUPAC | |
| 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 50 ng/ml |
| 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | — | 50 ng/ml |

MASS-LABELLED PCDDs & PCBs: SOLUTION/MIXTURES

Support solutions for **DFP-CVS-B10**, **DF-CVS-C10**, and **PCB-CVS-B10** (page 90)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------------------------|--|----------|
| DFP-IS-B10 | Mass-Labelled PCDD & PCB Internal Standard Solution | 1.2 ml |
| MASS-LABELLED PCDDs | | |
| | 1,3,7,8-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 ng/ml |
| | 1,2,4,7,8-Pentachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 ng/ml |
| | 1,2,3,4,6,8-Hexachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 ng/ml |
| | 1,2,3,4,6,7,9-Heptachloro[¹³ C ₁₂]dibenzo-p-dioxin | 10 ng/ml |
| MASS-LABELLED PCBs | | |
| | IUPAC | |
| | 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L |
| | 2,3,3',5,5-Pentachloro[¹³ C ₁₂]biphenyl | 111L |
| | 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L |
| | 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L |
| | | 10 ng/ml |
| | | 10 ng/ml |
| | | 10 ng/ml |
| | | 10 ng/ml |
| DFP-SS-A10 | Mass-Labelled PCDD & PCB Sampling Spike Solution | 1.2 ml |
| MASS-LABELLED PCDD & PCB | | |
| | IUPAC | |
| | 1,2,3,4-Tetrachloro[¹³ C ₁₂]dibenzo-p-dioxin | — |
| | 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L |
| | | 10 ng/ml |
| | | 10 ng/ml |

NATIVE PCDDs & PCDFs: SOLUTION/MIXTURES

(*) Support solutions for **DF-CVS-A10** and **DF-CVS-B10**

(**) Support solution for **DF-CVS-C10** and **DFP-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|-----------------------------------|----------|
| DF-ST-A* | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-ST-B* | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| DF-ST-C** | Native PCDD/PCDF Solution/Mixture | 1.2 ml |

| | DF-ST-A* | DF-ST-B* | DF-ST-C** |
|---|----------|----------|-----------|
| NATIVE PCDDs | (µg/ml) | (µg/ml) | (µg/ml) |
| 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 1.0 | 1.0 | 1.0 |
| 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 1.0 | 1.0 | 1.0 |
| 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 1.0 | 1.0 | — |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 1.0 | 1.0 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 1.0 | 1.0 | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 1.0 | 2.0 | 2.0 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 1.0 | 2.0 | 2.0 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 1.0 | 2.0 | 2.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 1.0 | 2.0 | 2.0 |
| Octachlorodibenzo-p-dioxin | 2.0 | 5.0 | 5.0 |
| NATIVE PCDFs | | | |
| 1,2,7,8-Tetrachlorodibenzofuran | 1.0 | 1.0 | — |
| 1,2,8,9-Tetrachlorodibenzofuran | 1.0 | 1.0 | 1.0 |
| 1,3,6,8-Tetrachlorodibenzofuran | 1.0 | 1.0 | 1.0 |
| 2,3,7,8-Tetrachlorodibenzofuran | 1.0 | 1.0 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 1.0 | 1.0 | 1.0 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 1.0 | 1.0 | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 1.0 | 2.0 | 2.0 |
| Octachlorodibenzofuran | 2.0 | 5.0 | 5.0 |

NATIVE PCDDs & PCDFs: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|-----------------------------------|----------|
| NK-ST-A | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| NK-ST-A4 | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| NK-ST-B2 | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| NK-ST-B4 | Native PCDD/PCDF Solution/Mixture | 1.2 ml |

| | NK-ST-A (µg/ml) | NK-ST-A4 (ng/ml) | NK-ST-B2 (ng/ml) | NK-ST-B4 (µg/ml) |
|---|--------------------|---------------------|---------------------|---------------------|
| NATIVE PCDDs | | | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 2.0 | 2.0 | 100 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 2.0 | 2.0 | 100 | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 2.0 | 2.0 | 200 | 2.0 |
| Octachlorodibenzo-p-dioxin | 4.0 | 4.0 | 400 | 5.0 |
| NATIVE PCDFs | | | | |
| 2,3,7,8-Tetrachlorodibenzofuran | 2.0 | 2.0 | 100 | 1.0 |
| 1,2,3,7,8-Pentachlorodibenzofuran | 2.0 | 2.0 | 100 | 1.0 |
| 2,3,4,7,8-Pentachlorodibenzofuran | 2.0 | 2.0 | 100 | 1.0 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 2.0 | 2.0 | 200 | 2.0 |
| Octachlorodibenzofuran | 4.0 | 4.0 | 400 | 5.0 |

PCDDs & PCDFs: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane/toluene solution) | Qty/Conc |
|------------------|--|----------|
| DDF-MDT | Native PCDD/PCDF Solution/Mixture | 1.2 ml |
| MDDF-MDT | Mass-Labelled PCDD/PCDF Solution/Mixture | 1.2 ml |

| NATIVE PCDDs & PCDFs | | DDF-MDT (µg/ml) |
|---------------------------------|--|---------------------------|
| Dibenzo-p-dioxin | | 1.0 |
| 2-Chlorodibenzo-p-dioxin | | 1.0 |
| 2,3-Dichlorodibenzo-p-dioxin | | 1.0 |
| 2,3,7-Trichlorodibenzo-p-dioxin | | 1.0 |
| Dibenzofuran | | 1.0 |
| 2-Chlorodibenzofuran | | 1.0 |
| 2,3-Dichlorodibenzofuran | | 1.0 |
| 2,3,8-Trichlorodibenzofuran | | 1.0 |

| MASS-LABELLED PCDDs & PCDFs | | MDDF-MDT (µg/ml) |
|--|--|----------------------------|
| [¹³ C ₁₂]Dibenzo-p-dioxin | | 1.0 |
| 2-Chloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1.0 |
| 2,3-Dichloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1.0 |
| 2,3,7-Trichloro[¹³ C ₁₂]dibenzo-p-dioxin | | 1.0 |
| [¹³ C ₁₂]Dibenzofuran | | 1.0 |
| 2-Chloro[¹³ C ₁₂]dibenzofuran | | 1.0 |
| 2,3-Dichloro[¹³ C ₁₂]dibenzofuran | | 1.0 |
| 2,3,8-Trichloro[¹³ C ₁₂]dibenzofuran | | 1.0 |

PCBs: ANALYTICAL METHOD SOLUTIONS

The electrophilic chlorination of biphenyl leads to complex mixtures of chlorobiphenyl (PCB) congeners ranging from monochloro- up to decachloro-biphenyl. Although 209 PCB congeners are theoretically possible, only about 150 are found in industrial mixtures and have ended up in the environment.

Certain PCB congeners are considered more toxic than others and are termed 'dioxin-like'. For this reason there are analytical methods that target only these 'dioxin-like' PCBs, while other methods focus on the major PCB congeners in the industrial mixtures. In addition, there are methods that address both groups of PCB congeners.

Wellington has designed, prepared, and offers a number of sets of calibration kits and support solutions for the analysis of individual PCB congeners.

WP-CVS

This set of solutions is to be used for the analysis of the 12 dioxin-like PCB congeners by HRGC/HRMS.

EPA Method 1668C

This series of calibration solutions, and corresponding support solutions, were prepared to be used according to U.S. EPA Method 1668, Revision C.

EPA Method 1668

This calibration kit and support solutions were designed and prepared to be used with the Draft version (March, 1997) of U.S. EPA Method 1668 which is still popular with some laboratories.

EC-9605-CVS

Environment Canada Method 1/RM/31 is a HRGC/LRMS method for PCB analysis and these solutions were prepared to be used with this method.

P48-W-CVS and P48-M-CVS

European Standard Method EN 1948-4 is to be used for analysis of the 12 'dioxin-like' PCB congeners and the 6 'marker' PCB congeners in stationary source emissions. These two calibration kits, and their support solutions, were prepared for this method.

WM48-CVS

This calibration set is a combination of P48-W-CVS and P48-M-CVS.



WP-CVS STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------------|
| WP-CVS | Dioxin-Like PCBs Calibration and Verification Solutions CS1-CS7 | 1 kit (7 ampoules) |
| WP-CS1 | CS1 | 500 µl |
| WP-CS2 | CS2 | 500 µl |
| WP-CS3 | CS3 | 500 µl |
| WP-CS4 | CS4 | 500 µl |
| WP-CS5 | CS5 | 500 µl |
| WP-CS6 | CS6 | 500 µl |
| WP-CS7 | CS7 | 500 µl |

| NATIVE PCBs | IUPAC | WP-CS1 | WP-CS2 | WP-CS3 | WP-CS4 | WP-CS5 | WP-CS6 | WP-CS7 |
|---|-------|---------|---------|---------|---------|---------|---------|---------|
| | | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 | 800 |
| MASS-LABELLED PCBs | | | | | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| INTERNAL STANDARDS: MASS-LABELLED PCBs | | | | | | | | |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 50 | 50 | 50 | 50 | 50 | 50 | 50 |

WP-CVS STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|----------------------------|----------|
| WP-LCS | Surrogate Spiking Solution | 1.2 ml |
| WP-ISS | Internal Standard Solution | 1.2 ml |
| WP-STK | Native PCB Solution | 1.2 ml |

| NATIVE PCBs | IUPAC | WP-LCS (ng/ml) | WP-ISS (ng/ml) | WP-STK (ng/ml) |
|---|-------|-------------------|-------------------|-------------------|
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | — | — | 2000 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | — | — | 2000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | — | — | 2000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | — | — | 2000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | — | — | 2000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | — | — | 2000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | — | — | 2000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | — | — | 2000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | — | — | 2000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | — | — | 2000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | — | — | 2000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | — | — | 2000 |
| MASS-LABELLED PCBs | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 1000 | — | — |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 1000 | — | — |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 1000 | — | — |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 1000 | — | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 1000 | — | — |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 1000 | — | — |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 1000 | — | — |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 1000 | — | — |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 1000 | — | — |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 1000 | — | — |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 1000 | — | — |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 1000 | — | — |
| INTERNAL STANDARDS: MASS-LABELLED PCBs | | | | |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | — | 1000 | — |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | — | 1000 | — |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | — | 1000 | — |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | — | 1000 | — |

EPA METHOD 1668C STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------------|
| 68C-CV5 | EPA Method 1668C Calibration and Verification Solutions CS0.2-CS5 | 1 kit (6 ampoules) |
| 68C-CS0.2 | CS0.2 High Sensitivity | 200 µl |
| 68C-CS1 | CS1 | 200 µl |
| 68C-CS2 | CS2 | 200 µl |
| 68C-CS3 | CS3 Calibration Verification | 500 µl |
| 68C-CS4 | CS4 | 200 µl |
| 68C-CS5 | CS5 | 200 µl |

NOTE: The above product codes were updated to reflect the change of EPA Method 1668B to 1668C in April of 2010.

| Native Toxics/LOC | IUPAC | 68C-CS0.2 (ng/ml) | 68C-CS1 (ng/ml) | 68C-CS2 (ng/ml) | 68C-CS3 (ng/ml) | 68C-CS4 (ng/ml) | 68C-CS5 (ng/ml) |
|---|-------|----------------------|--------------------|--------------------|--------------------|--------------------|--------------------|
| 2-Chlorobiphenyl | 1 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 4-Chlorobiphenyl | 3 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2'-Dichlorobiphenyl | 4 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 4,4'-Dichlorobiphenyl | 15 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',6-Trichlorobiphenyl | 19 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,4,4'-Trichlorobiphenyl | 37 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',6,6'-Tetrachlorobiphenyl | 54 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',4,6,6'-Pentachlorobiphenyl | 104 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',4,4',6,6'-Hexachlorobiphenyl | 155 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 188 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 202 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5,5',6-Octachlorobiphenyl | 205 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 208 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| Decachlorobiphenyl | 209 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| Labelled Toxics/LOC/Window-Defining (68C-LCS) | | | | | | | |
| 2-Chloro[¹³ C ₁₂]biphenyl | 1L | 100 | 100 | 100 | 100 | 100 | 100 |
| 4-Chloro[¹³ C ₁₂]biphenyl | 3L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2'-Dichloro[¹³ C ₁₂]biphenyl | 4L | 100 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-Dichloro[¹³ C ₁₂]biphenyl | 15L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',6-Trichloro[¹³ C ₁₂]biphenyl | 19L | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 37L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',6,6'-Tetrachloro[¹³ C ₁₂]biphenyl | 54L | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,6,6'-Pentachloro[¹³ C ₁₂]biphenyl | 104L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',6,6'-Hexachloro[¹³ C ₁₂]biphenyl | 155L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4',5,6,6'-Heptachloro[¹³ C ₁₂]biphenyl | 188L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',5,5',6,6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5',6-Octachloro[¹³ C ₁₂]biphenyl | 205L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,5,5',6,6'-Nonachloro[¹³ C ₁₂]biphenyl | 208L | 100 | 100 | 100 | 100 | 100 | 100 |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 100 | 100 | 100 | 100 | 100 | 100 |
| Labelled Clean-Up (68C-CS) | | | | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | 100 | 100 | 100 | 100 | 100 | 100 |
| Labelled Injection/Internal (68C-IS) | | | | | | | |
| 2,5-Dichloro[¹³ C ₁₂]biphenyl | 9L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | 100 | 100 | 100 | 100 | 100 | 100 |

EPA METHOD 1668C STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| 68C-LCS | Labelled Toxics/LOC/Window Defining Stock Solution | 1.2 ml |
| 68C-CS | Labelled Cleanup Stock Solution | 1.2 ml |
| 68C-IS | Labelled Injection/Internal Standard Stock Solution | 1.2 ml |
| 68C-PAR | Native Toxics/LOC Stock Solution | 1.2 ml |

NOTE: The above product codes were updated to reflect the change of EPA Method 1668B to 1668C in April of 2010.

| Native Toxics/LOC | IUPAC | 68C-LCS (ng/ml) | 68C-CS (ng/ml) | 68C-IS (ng/ml) | 68C-PAR (ng/ml) |
|---|-------|--------------------|-------------------|-------------------|--------------------|
| 2-Chlorobiphenyl | 1 | — | — | — | 2000 |
| 4-Chlorobiphenyl | 3 | — | — | — | 2000 |
| 2,2'-Dichlorobiphenyl | 4 | — | — | — | 2000 |
| 4,4'-Dichlorobiphenyl | 15 | — | — | — | 2000 |
| 2,2',6-Trichlorobiphenyl | 19 | — | — | — | 2000 |
| 3,4,4'-Trichlorobiphenyl | 37 | — | — | — | 2000 |
| 2,2',6,6'-Tetrachlorobiphenyl | 54 | — | — | — | 2000 |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | — | — | — | 2000 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | — | — | — | 2000 |
| 2,2',4,6,6'-Pentachlorobiphenyl | 104 | — | — | — | 2000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | — | — | — | 2000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | — | — | — | 2000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | — | — | — | 2000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | — | — | — | 2000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | — | — | — | 2000 |
| 2,2',4,4',6,6'-Hexachlorobiphenyl | 155 | — | — | — | 2000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | — | — | — | 2000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | — | — | — | 2000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | — | — | — | 2000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | — | — | — | 2000 |
| 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 188 | — | — | — | 2000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | — | — | — | 2000 |
| 2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 202 | — | — | — | 2000 |
| 2,3,3',4,4',5,5',6-Octachlorobiphenyl | 205 | — | — | — | 2000 |
| 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | — | — | — | 2000 |
| 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 208 | — | — | — | 2000 |
| Decachlorobiphenyl | 209 | — | — | — | 2000 |
| Labelled Toxics/LOC/Window-Defining (68C-LCS) | | | | | |
| 2-Chloro[¹³ C ₁₂]biphenyl | 1L | 1000 | — | — | — |
| 4-Chloro[¹³ C ₁₂]biphenyl | 3L | 1000 | — | — | — |
| 2,2'-Dichloro[¹³ C ₁₂]biphenyl | 4L | 1000 | — | — | — |
| 4,4'-Dichloro[¹³ C ₁₂]biphenyl | 15L | 1000 | — | — | — |
| 2,2',6-Trichloro[¹³ C ₁₂]biphenyl | 19L | 1000 | — | — | — |
| 3,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 37L | 1000 | — | — | — |
| 2,2',6,6'-Tetrachloro[¹³ C ₁₂]biphenyl | 54L | 1000 | — | — | — |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 1000 | — | — | — |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 1000 | — | — | — |
| 2,2',4,6,6'-Pentachloro[¹³ C ₁₂]biphenyl | 104L | 1000 | — | — | — |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 1000 | — | — | — |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 1000 | — | — | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 1000 | — | — | — |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 1000 | — | — | — |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 1000 | — | — | — |
| 2,2',4,4',6,6'-Hexachloro[¹³ C ₁₂]biphenyl | 155L | 1000 | — | — | — |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 1000 | — | — | — |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 1000 | — | — | — |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 1000 | — | — | — |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 1000 | — | — | — |
| 2,2',3,4',5,6,6'-Heptachloro[¹³ C ₁₂]biphenyl | 188L | 1000 | — | — | — |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 1000 | — | — | — |
| 2,2',3,3',5,5',6,6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | 1000 | — | — | — |
| 2,3,3',4,4',5,5',6-Octachloro[¹³ C ₁₂]biphenyl | 205L | 1000 | — | — | — |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | 1000 | — | — | — |
| 2,2',3,3',4,5,5',6,6'-Nonachloro[¹³ C ₁₂]biphenyl | 208L | 1000 | — | — | — |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 1000 | — | — | — |
| Labelled Clean-Up (68C-CS) | | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | — | 1000 | — | — |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | — | 1000 | — | — |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | — | 1000 | — | — |
| Labelled Injection/Internal (68C-IS) | | | | | |
| 2,5-Dichloro[¹³ C ₁₂]biphenyl | 9L | — | — | 5000 | — |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | — | — | 5000 | — |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | — | — | 5000 | — |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | — | — | 5000 | — |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | — | — | 5000 | — |

EPA METHOD 1668 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|---|-----------------------|
| EPA-1668CVS | EPA Method 1668 Calibration and Verification Solutions CS1-CS5 | 1 kit (5 ampoules) |
| EPA-1668CS1 | CS1 | 200 µl |
| EPA-1668CS2 | CS2 | 200 µl |
| EPA-1668CS3 | CS3 Calibration Verification | 500 µl |
| EPA-1668CS4 | CS4 | 200 µl |
| EPA-1668CS5 | CS5 | 200 µl |

| NATIVE PCBs | IUPAC | 1668CS1 (ng/ml) | 1668CS2 (ng/ml) | 1668CS3 (ng/ml) | 1668CS4 (ng/ml) | 1668CS5 (ng/ml) |
|---|-------|--------------------|--------------------|--------------------|--------------------|--------------------|
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 2.5 | 10 | 50 | 200 | 1000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 2.5 | 10 | 50 | 200 | 1000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 5.0 | 20 | 100 | 400 | 2000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 5.0 | 20 | 100 | 400 | 2000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 5.0 | 20 | 100 | 400 | 2000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 5.0 | 20 | 100 | 400 | 2000 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 5.0 | 20 | 100 | 400 | 2000 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 5.0 | 20 | 100 | 400 | 2000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 5.0 | 20 | 100 | 400 | 2000 |
| MASS-LABELLED PCBs | | | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 100 | 100 | 100 | 100 | 100 |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 200 | 200 | 200 | 200 | 200 |
| CLEANUP STANDARDS | | | | | | |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 2.5 | 10 | 50 | 200 | 1000 |
| INTERNAL STANDARDS | | | | | | |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | 100 | 100 | 100 | 100 | 100 |

EPA METHOD 1668 STANDARD SOLUTIONS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|----------------------------------|----------|
| EPA-1668LCS | Labelled Compound Stock Solution | 1.2 ml |
| EPA-1668CS | Cleanup Standard Solution | 1.2 ml |
| EPA-1668IS | Internal Standard Stock Solution | 1.2 ml |
| EPA-1668PAR | Precision and Recovery Solution | 1.2 ml |

| NATIVE PCBs | IUPAC | 1668LCS (ng/ml) | 1668CS (ng/ml) | 1668IS (ng/ml) | 1668PAR (ng/ml) |
|---|-------|--------------------|-------------------|-------------------|--------------------|
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | — | — | — | 20 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | — | — | — | 1000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | — | — | — | 1000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | — | — | — | 1000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | — | — | — | 1000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | — | — | — | 100 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | — | — | — | 1000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | — | — | — | 1000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | — | — | — | 1000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | — | — | — | 200 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | — | — | — | 200 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | — | — | — | 1000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | — | — | — | 200 |
| MASS-LABELLED PCBs | | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 1000 | — | — | — |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 1000 | — | — | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 1000 | — | — | — |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 1000 | — | — | — |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 1000 | — | — | — |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 1000 | — | — | — |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 1000 | — | — | — |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 1000 | — | — | — |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 1000 | — | — | — |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 1000 | — | — | — |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 2000 | — | — | — |
| CLEANUP STANDARDS | | | | | |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | — | 200 | — | — |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | — | 1000 | — | — |
| INTERNAL STANDARDS | | | | | |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | — | — | 1000 | — |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | — | — | 1000 | — |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | — | — | 1000 | — |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | — | — | 1000 | — |

ENVIRONMENT CANADA METHOD 1/RM/31 STANDARD SOLUTIONS

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|-------------------|---|-----------------------|
| EC9605-CVS | PCB Calibration Solutions for GC/MS Calibration and Verification Solutions CS1-CS5 | 1 kit (5 ampoules) |
| ECPCS1 | CS1 | 500 µl |
| ECPCS2 | CS2 | 500 µl |
| ECPCS3 | CS3 | 500 µl |
| ECPCS4 | CS4 | 500 µl |
| ECPCS5 | CS5 | 500 µl |

| NATIVE PCBs* | IUPAC | ECPCS1 (ng/ml) | ECPCS2 (ng/ml) | ECPCS3 (ng/ml) | ECPCS4 (ng/ml) | ECPCS5 (ng/ml) |
|---|--------------|---------------------------|---------------------------|---------------------------|---------------------------|---------------------------|
| 2,2',5-Trichlorobiphenyl | 18 | 20 | 50 | 200 | 800 | 2000 |
| 2,4,4'-Trichlorobiphenyl | 28 | 20 | 50 | 200 | 800 | 2000 |
| 2',3,4-Trichlorobiphenyl | 33 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',5,5'-Tetrachlorobiphenyl | 52 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,5'-Tetrachlorobiphenyl | 44 | 20 | 50 | 200 | 800 | 2000 |
| 2,3',4',5-Tetrachlorobiphenyl | 70 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',4,5,5'-Pentachlorobiphenyl | 101 | 20 | 50 | 200 | 800 | 2000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 20 | 50 | 200 | 800 | 2000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,4'-Hexachlorobiphenyl | 128 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,4',5,5',6-Heptachlorobiphenyl | 187 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,5,5',6'-Octachlorobiphenyl | 199 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,4',5,6-Octachlorobiphenyl | 195 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,4',5,5'-Octachlorobiphenyl | 194 | 20 | 50 | 200 | 800 | 2000 |
| 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | 20 | 50 | 200 | 800 | 2000 |
| Decachlorobiphenyl | 209 | 20 | 50 | 200 | 800 | 2000 |
| MASS-LABELLED PCBs* | | | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 400 | 400 | 400 | 400 | 400 |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 400 | 400 | 400 | 400 | 400 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 400 | 400 | 400 | 400 | 400 |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 400 | 400 | 400 | 400 | 400 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 400 | 400 | 400 | 400 | 400 |
| 2,2',3,3',5,5',6,6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | 400 | 400 | 400 | 400 | 400 |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | 400 | 400 | 400 | 400 | 400 |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 400 | 400 | 400 | 400 | 400 |
| MASS-LABELLED RECOVERY STANDARDS* | | | | | | |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 400 | 400 | 400 | 400 | 400 |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | 400 | 400 | 400 | 400 | 400 |

* In order of elution on a 60m DB-5 column.

ENVIRONMENT CANADA METHOD 1/RM/31 STANDARD SOLUTIONS

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|-------------------|---------------------------------|----------|
| EC9605-RS | Recovery Standard Solution | 1.2 ml |
| EC9605-SS | Surrogate Solution | 1.2 ml |
| EC9605-PAR | Precision and Recovery Solution | 1.2 ml |

| NATIVE PCBs | IUPAC | EC9605-RS (µg/ml) | EC9605-SS (µg/ml) | EC9605-PAR (ng/ml) |
|---|-------|----------------------|----------------------|-----------------------|
| 2,2',5-Trichlorobiphenyl | 18 | — | — | 100 |
| 2,4,4'-Trichlorobiphenyl | 28 | — | — | 100 |
| 2',3,4-Trichlorobiphenyl | 33 | — | — | 100 |
| 2,2',5,5'-Tetrachlorobiphenyl | 52 | — | — | 100 |
| 2,2',3,5'-Tetrachlorobiphenyl | 44 | — | — | 100 |
| 2,3',4',5-Tetrachlorobiphenyl | 70 | — | — | 100 |
| 2,2',4,5,5'-Pentachlorobiphenyl | 101 | — | — | 100 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | — | — | 100 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | — | — | 100 |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | — | — | 100 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | — | — | 100 |
| 2,2',3,3',4,4'-Hexachlorobiphenyl | 128 | — | — | 100 |
| 2,2',3,4',5,5',6-Heptachlorobiphenyl | 187 | — | — | 100 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | — | — | 100 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | — | — | 100 |
| 2,2',3,3',4,5,5',6'-Octachlorobiphenyl | 199 | — | — | 100 |
| 2,2',3,3',4,4',5,6-Octachlorobiphenyl | 195 | — | — | 100 |
| 2,2',3,3',4,4',5,5'-Octachlorobiphenyl | 194 | — | — | 100 |
| 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | — | — | 100 |
| Decachlorobiphenyl | 209 | — | — | 100 |
| MASS-LABELLED PCBs | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | — | 2.0 | — |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | — | 2.0 | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | — | 2.0 | — |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | — | 2.0 | — |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | — | 2.0 | — |
| 2,2',3,3',5,5',6,6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | — | 2.0 | — |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | — | 2.0 | — |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | — | 2.0 | — |
| MASS-LABELLED RECOVERY STANDARDS | | | | |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 2.0 | — | — |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | 2.0 | — | — |

P48-W-CVS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-----------------------|
| P48-W-CVS | P48-W-CVS; EN 1948-4:2010 HRGC/HRMS Calibration Solutions for the Dioxin-like PCBs | 1 kit (6 ampoules) |
| P48-W-CS1 | CS1 | 500 µl |
| P48-W-CS2 | CS2 | 500 µl |
| P48-W-CS3 | CS3 | 500 µl |
| P48-W-CS4 | CS4 | 500 µl |
| P48-W-CS5 | CS5 | 500 µl |
| P48-W-CS6 | CS6 | 500 µl |

| NATIVE DIOXIN-LIKE PCB CONGENERS | IUPAC | P48-W-CS1 (pg/µl) | P48-W-CS2 (pg/µl) | P48-W-CS3 (pg/µl) | P48-W-CS4 (pg/µl) | P48-W-CS5 (pg/µl) | P48-W-CS6 (pg/µl) |
|--------------------------------------|-------|----------------------|----------------------|----------------------|----------------------|----------------------|----------------------|
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.6 | 6.0 | 60 | 300 | 1200 | 4800 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.1 | 1.0 | 10 | 50 | 200 | 800 |

WHO PCB EXTRACTION SPIKE (P48-W-ES)

| | | | | | | | |
|---|------|----|----|----|----|----|----|
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 10 | 10 | 10 | 10 | 10 | 10 |

SAMPLING SPIKE (P48-SS)

| | | | | | | | |
|---|------|----|----|----|----|----|----|
| 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 60L | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 127L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 159L | 10 | 10 | 10 | 10 | 10 | 10 |

RECOVERY SPIKE (P48-RS)

| | | | | | | | |
|---|------|----|----|----|----|----|----|
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 | 10 | 10 | 10 | 10 | 10 |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|--|-----------------------|
| P48-M-CVS | P48-M-CVS; EN 1948-4:2010 HRGC/HRMS Calibration Solutions for the Marker PCBs | 1 kit (6 ampoules) |
| P48-M-CS0.1 | CS0.1 | 500 µl |
| P48-M-CS1 | CS1 | 500 µl |
| P48-M-CS2 | CS2 | 500 µl |
| P48-M-CS3 | CS3 | 500 µl |
| P48-M-CS4 | CS4 | 500 µl |
| P48-M-CS5 | CS5 | 500 µl |

| NATIVE MARKER PCB CONGENERS | IUPAC | P48-M-CS0.1 (pg/µl) | P48-M-CS1 (pg/µl) | P48-M-CS2 (pg/µl) | P48-M-CS3 (pg/µl) | P48-M-CS4 (pg/µl) | P48-M-CS5 (pg/µl) |
|---|-------|---------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| 2,4,4'-Trichlorobiphenyl | 28 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| 2,2',5,5'-Tetrachlorobiphenyl | 52 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| 2,2',4,5,5'-Pentachlorobiphenyl | 101 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 0.1 | 1.0 | 10 | 100 | 500 | 5000 |
| MARKER PCB EXTRACTION SPIKE (P48-M-ES) | | | | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 100 | 100 | 100 | 100 | 100 | 100 |
| SAMPLING SPIKE (P48-SS) | | | | | | | |
| 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 60L | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 127L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 159L | 10 | 10 | 10 | 10 | 10 | 10 |
| RECOVERY SPIKE (P48-RS) | | | | | | | |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 | 10 | 10 | 10 | 10 | 10 |

WM48-CVS

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------------|
| WM48-CVS | WM48-CVS; EN 1948-4:2010 HRGC/HRMS Calibration Solutions for the Dioxin-like and Marker PCBs | 1 kit (6 ampoules) |
| WM48-CS1 | CS1 | 500 µl |
| WM48-CS2 | CS2 | 500 µl |
| WM48-CS3 | CS3 | 500 µl |
| WM48-CS4 | CS4 | 500 µl |
| WM48-CS5 | CS5 | 500 µl |
| WM48-CS6 | CS6 | 500 µl |

| | | WM48- CS1 (pg/µl) | WM48- CS2 (pg/µl) | WM48- CS3 (pg/µl) | WM48- CS4 (pg/µl) | WM48- CS5 (pg/µl) | WM48- CS6 (pg/µl) |
|---|---|-------------------------|-------------------------|-------------------------|-------------------------|-------------------------|-------------------------|
| NATIVE DIOXIN-LIKE PCB CONGENERS | IUPAC | | | | | | |
| | 3,3',4,4'-Tetrachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 3,4,4',5-Tetrachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3,3',4,4'-Pentachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3,4,4',5-Pentachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3',4,4',5-Pentachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2',3,4,4',5-Pentachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 3,3',4,4',5-Pentachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3,3',4,4',5-Hexachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3,3',4,4',5'-Hexachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3',4,4',5,5'-Hexachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 3,3',4,4',5,5'-Hexachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| | 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 0.1 | 0.5 | 2 | 10 | 40 | 200 |
| NATIVE MARKER PCB CONGENERS | | | | | | | |
| | 2,4,4'-Trichlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2,2',5,5'-Tetrachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2,2',4,5,5'-Pentachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2,2',3,4,4',5'-Hexachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2,2',4,4',5,5'-Hexachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| | 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 0.5 | 2.5 | 10 | 50 | 200 | 1000 |
| EXTRACTION SPIKE | | | | | | | |
| | 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| SAMPLING SPIKE | | | | | | | |
| | 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 3,3',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| RECOVERY SPIKE | | | | | | | |
| | 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |
| | 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 10 | 10 | 10 | 10 | 10 | 10 |

P48-W-PAR: Native Dioxin-Like (WHO) PCB Solution

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|---------------------------|--------------------------|
| P48-W-PAR | P48-W-PAR; EN 1948-4:2010 | 1.2 ml |
| NATIVE DIOXIN-LIKE (WHO) PCB CONGENERS | IUPAC | P48-W-PAR (pg/μl) |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 250 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 250 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 250 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 250 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 1500 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 250 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 250 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 250 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 250 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 250 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 250 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 250 |

P48-M-PAR: Native Marker PCB Solution

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------------------------|---------------------------|--------------------------|
| P48-M-PAR | P48-M-PAR; EN 1948-4:2010 | 1.2 ml |
| NATIVE MARKER PCB CONGENERS | IUPAC | P48-M-PAR (pg/μl) |
| 2,4,4'-Trichlorobiphenyl | 28 | 250 |
| 2,2',5,5'-Tetrachlorobiphenyl | 52 | 250 |
| 2,2',4,5,5'-Pentachlorobiphenyl | 101 | 250 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | 250 |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | 250 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 250 |

SUPPORT SOLUTIONS FOR EN 1948-4:2010

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| P48-W-ES | Dioxin-like (WHO) PCB Extraction Standard | 1.2 ml |
| P48-M-ES | Marker PCB Extraction Standard | 1.2 ml |
| P48-SS | Mass-Labelled PCB Sampling Standard | 1.2 ml |
| P48-RS | Mass-Labelled PCB Recovery Standard | 1.2 ml |

| DIOXIN-LIKE PCB EXTRACTION STANDARD | IUPAC | P48-W-ES (pg/µl) | P48-M-ES (pg/µl) | P48-SS (pg/µl) | P48-RS (pg/ul) |
|---|-------|---------------------|---------------------|-------------------|-------------------|
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 100 | — | — | — |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 100 | — | — | — |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 100 | — | — | — |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 100 | — | — | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 100 | — | — | — |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 100 | — | — | — |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 100 | — | — | — |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 100 | — | — | — |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 100 | — | — | — |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 100 | — | — | — |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 100 | — | — | — |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 100 | — | — | — |
| MARKER PCB EXTRACTION STANDARD | | | | | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | — | 1000 | — | — |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | — | 1000 | — | — |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | — | 1000 | — | — |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | — | 1000 | — | — |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | — | 1000 | — | — |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | — | 1000 | — | — |
| MASS-LABELLED PCB SAMPLING STANDARD | | | | | |
| 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 60L | — | — | 100 | — |
| 3,3',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 127L | — | — | 100 | — |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 159L | — | — | 100 | — |
| MASS-LABELLED PCB RECOVERY STANDARD | | | | | |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | — | — | — | 100 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | — | — | — | 100 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | — | — | — | 100 |

PCBs: MASS-LABELLED CONGENERES

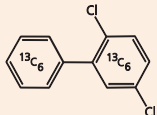
Wellington has prepared a very comprehensive collection of fully labelled ($^{13}\text{C}_{12}$) individual PCB congeners including the 12 'dioxin-like' PCBs and the more prominent congeners found in commercial mixtures and the environment.

All of the ^{13}C -PCBs in the following pages were prepared using one-product, unambiguous routes and purified using a variety of methods. Their structures and chemical and isotopic purities were confirmed using various instruments and this data is included in the Certificates of Analysis.

Additional ^{13}C -PCBs may be added in the future, so please continue to visit our website for updates or contact **Wellington** or your local distributor if you have any specific requests.



MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|---|
| MBP-1 |  <p>2-Chloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-3 |  <p>4-Chloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-4 |  <p>2,2'-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-5 |  <p>2,3-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-8 |  <p>2,4'-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-9 |  <p>2,5-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-11 |  <p>3,3'-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-15 |  <p>4,4'-Dichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-19 |  <p>2,2',6-Trichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-28 |  <p>2,4,4'-Trichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

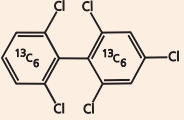
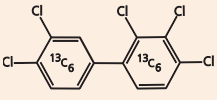
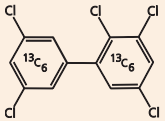
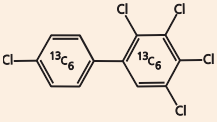
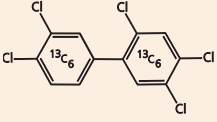
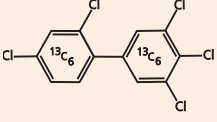
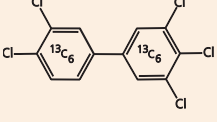
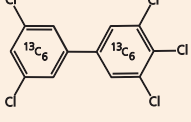
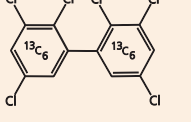
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|--|
| MBP-31 |  <p>2,4',5-Trichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-37 |  <p>3,4,4'-Trichloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-52 |  <p>2,2',5,5'-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-54 |  <p>2,2',6,6'-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-60 |  <p>2,3,4,4'-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-70 |  <p>2,3',4',5-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-77 |  <p>3,3',4,4'-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-79 |  <p>3,3',4,5'-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-81 |  <p>3,4,4',5-Tetrachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-95 |  <p>2,2',3,5',6-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|--|
| MBP-101 |  <p>2,2',4,5,5'-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-104 |  <p>2,2',4,6,6'-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-105 |  <p>2,3,3',4,4'-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-111 |  <p>2,3,3',5,5'-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-114 |  <p>2,3,4,4',5-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-118 |  <p>2,3',4,4',5-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-123 |  <p>2',3,4,4',5-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-126 |  <p>3,3',4,4',5-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-127 |  <p>3,3',4,5,5'-Pentachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-133 |  <p>2,2',3,3',5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

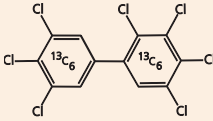
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|---|
| MBP-138 |  <p>2,2',3,4,4',5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-153 |  <p>2,2',4,4',5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-155 |  <p>2,2',4,4',6,6'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-156 |  <p>2,3,3',4,4',5-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-157 |  <p>2,3,3',4,4',5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-159 |  <p>2,3,3',4,5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-162 |  <p>2,3,3',4',5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-167 |  <p>2,3',4,4',5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-169 |  <p>3,3',4,4',5,5'-Hexachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-170 |  <p>2,2',3,3',4,4',5-Heptachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|---|
| MBP-178 |  <p>2,2',3,3',5,5',6-Heptachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-180 |  <p>2,2',3,4,4',5,5'-Heptachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-188 |  <p>2,2',3,4',5,6,6'-Heptachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-189 |  <p>2,3,3',4,4',5,5'-Heptachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-194 |  <p>2,2',3,3',4,4',5,5'-Octachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-202 |  <p>2,2',3,3',5,5',6,6'-Octachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-205 |  <p>2,3,3',4,4',5,5',6-Octachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-206 |  <p>2,2',3,3',4,4',5,5',6-Nonachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-208 |  <p>2,2',3,3',4,5,5',6,6'-Nonachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBP-209 |  <p>Decachloro[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

PCBs: SPECIALTY SOLUTION/MIXTURES

A number of additional PCB calibration sets and other solution/mixtures have been prepared and are presented in this section.

PCB-CVS-A10 and PCB-CVS-B10, and their support solutions, were designed and prepared to be used to satisfy the requirements of the Japanese Industrial Standards **JIS K 0311:2005** and **JIS K 0312:2005**. Note that PCB congeners 170 and 180, and their ¹³C analogues, have been added to the 12 dioxin-like PCBs.

Calibration sets and mixes containing a larger number of PCB congeners, such as PCB-CVS-H, are also offered.

All of the solutions from this section are accompanied by detailed CofAs that include HRGC/LRMS and/or HRGC/HRMS data as appropriate, along with RRF summaries for the calibration sets.



PCB-CVS-H

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------------------|
| PCB-CVS-H | Calibration Solutions for HRGC/HRMS Analysis of Polychlorinated Biphenyls (PCBs) | 1 kit (6 x 200 µl ampoules) |
| PCB-CS1-H | CS1 | 200 µl |
| PCB-CS2-H | CS2 | 200 µl |
| PCB-CS3-H | CS3 | 200 µl |
| PCB-CS4-H | CS4 | 200 µl |
| PCB-CS5-H | CS5 | 200 µl |
| PCB-CS6-H | CS6 | 200 µl |

| | PCB-CS1-H (ng/ml) | PCB-CS2-H (ng/ml) | PCB-CS3-H (ng/ml) | PCB-CS4-H (ng/ml) | PCB-CS5-H (ng/ml) | PCB-CS6-H (ng/ml) |
|--|-------------------|-------------------|-------------------|-------------------|-------------------|-------------------|
| NATIVE CHLORINATED BIPHENYLS (IUPAC) | | | | | | |
| CHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 1, 3 | | | | | | |
| DICHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 4, 6, 8, 10, 15 | | | | | | |
| TRICHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 16, 18, 19, 22, 28, 31, 33, 37 | | | | | | |
| TETRACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 40, 41, 44, 49, 52, 54, 60, 66, 70, 74, 77, 81 | | | | | | |
| PENTACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 84, 85, 87, 90, 95, 97, 99, 101, 104, 105, 110, 114, 118, 119, 123, 126 | | | | | | |
| HEXACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 128, 129, 135, 137, 138, 141, 149, 151, 153, 155, 156, 157, 158, 167, 168, 169 | | | | | | |
| HEPTACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 170, 171, 174, 177, 178, 180, 183, 187, 188, 189, 191, 193 | | | | | | |
| OCTACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 194, 199, 200, 201, 202, 203, 205 | | | | | | |
| NONACHLOROBIPHENYLS | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 206, 207, 208 | | | | | | |
| DECACHLOROBIPHENYL | 0.1 | 0.5 | 2.0 | 10 | 40 | 200 |
| 209 | | | | | | |
| MASS-LABELLED CHLORINATED BIPHENYLS | IUPAC | | | | | |
| EXTRACTION STANDARDS | | | | | | |
| 2-Chloro[¹³ C ₁₂]biphenyl | 1L | 50 | 50 | 50 | 50 | 50 |
| 4-Chloro[¹³ C ₁₂]biphenyl | 3L | 50 | 50 | 50 | 50 | 50 |
| 2,2'-Dichloro[¹³ C ₁₂]biphenyl | 4L | 50 | 50 | 50 | 50 | 50 |
| 2,4'-Dichloro[¹³ C ₁₂]biphenyl | 8L | 50 | 50 | 50 | 50 | 50 |
| 4,4'-Dichloro[¹³ C ₁₂]biphenyl | 15L | 50 | 50 | 50 | 50 | 50 |
| 2,2',6-Trichloro[¹³ C ₁₂]biphenyl | 19L | 50 | 50 | 50 | 50 | 50 |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 50 | 50 | 50 | 50 | 50 |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 50 | 50 | 50 | 50 | 50 |
| 2,2',6,6'-Tetrachloro[¹³ C ₁₂]biphenyl | 54L | 50 | 50 | 50 | 50 | 50 |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 50 | 50 | 50 | 50 | 50 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,5',6-Pentachloro[¹³ C ₁₂]biphenyl | 95L | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,6,6'-Pentachloro[¹³ C ₁₂]biphenyl | 104L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 50 | 50 | 50 | 50 | 50 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 50 | 50 | 50 | 50 | 50 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 50 | 50 | 50 | 50 | 50 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',6,6'-Hexachloro[¹³ C ₁₂]biphenyl | 155L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 50 | 50 | 50 | 50 | 50 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4',5,6,6'-Heptachloro[¹³ C ₁₂]biphenyl | 188L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',5,5',6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,4',5,5',6-Octachloro[¹³ C ₁₂]biphenyl | 205L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',4,4',5,5',6,6'-Nonachloro[¹³ C ₁₂]biphenyl | 208L | 50 | 50 | 50 | 50 | 50 |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 50 | 50 | 50 | 50 | 50 |
| RECOVERY/INTERNAL STANDARDS | | | | | | |
| 2,5-Dichloro[¹³ C ₁₂]biphenyl | 9L | 50 | 50 | 50 | 50 | 50 |
| 3,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 37L | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 162L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | 50 | 50 | 50 | 50 | 50 |
| SAMPLING/CLEANUP STANDARDS | | | | | | |
| 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 60L | 50 | 50 | 50 | 50 | 50 |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 159L | 50 | 50 | 50 | 50 | 50 |

SUPPORT SOLUTIONS FOR PCB-CVS-H

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| PCB-LCS-H | Mass-Labelled PCB Extraction Standards | 1.2 ml |
| PCB-ISS-H | Mass-Labelled PCB Internal/Recovery Standards | 1.2 ml |
| PCB-SCS-H | Mass-Labelled PCB Cleanup/Sampling Standards | 1.2 ml |
| PCB-PAR-H | Native PCB Solution | 1.2 ml |

| NATIVE CHLORINATED BIPHENYLS (IUPAC) | PCB-LCS-H (ng/ml) | PCB-ISS-H (ng/ml) | PCB-SCS-H (ng/ml) | PCB-PAR-H (ng/ml) |
|--|-------------------|-------------------|-------------------|-------------------|
| CHLOROBIPHENYLS | — | — | — | 500 |
| 1, 3 | | | | |
| DICHLOROBIPHENYLS | — | — | — | 500 |
| 4, 6, 8, 10, 15 | | | | |
| TRICHLOROBIPHENYLS | — | — | — | 500 |
| 16, 18, 19, 22, 28, 31, 33, 37 | | | | |
| TETRACHLOROBIPHENYLS | — | — | — | 500 |
| 40, 41, 44, 49, 52, 54, 60, 66, 70, 74, 77, 81 | | | | |
| PENTACHLOROBIPHENYLS | — | — | — | 500 |
| 84, 85, 87, 90, 95, 97, 99, 101, 104, 105, 110, 114, 118, 119, 123, 126 | | | | |
| HEXACHLOROBIPHENYLS | — | — | — | 500 |
| 128, 129, 135, 137, 138, 141, 149, 151, 153, 155, 156, 157, 158, 167, 168, 169 | | | | |
| HEPTACHLOROBIPHENYLS | — | — | — | 500 |
| 170, 171, 174, 177, 178, 180, 183, 187, 188, 189, 191, 193 | | | | |
| OCTACHLOROBIPHENYLS | — | — | — | 500 |
| 194, 199, 200, 201, 202, 203, 205 | | | | |
| NONACHLOROBIPHENYLS | — | — | — | 500 |
| 206, 207, 208 | | | | |
| DECACHLOROBIPHENYL | — | — | — | 500 |
| 209 | | | | |
| MASS-LABELLED CHLORINATED BIPHENYLS EXTRACTION STANDARDS | IUPAC | | | |
| 2-Chloro[¹³ C ₁₂]biphenyl | 1L | 1000 | — | — |
| 4-Chloro[¹³ C ₁₂]biphenyl | 3L | 1000 | — | — |
| 2,2'-Dichloro[¹³ C ₁₂]biphenyl | 4L | 1000 | — | — |
| 2,4'-Dichloro[¹³ C ₁₂]biphenyl | 8L | 1000 | — | — |
| 4,4'-Dichloro[¹³ C ₁₂]biphenyl | 15L | 1000 | — | — |
| 2,2',6-Trichloro[¹³ C ₁₂]biphenyl | 19L | 1000 | — | — |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 1000 | — | — |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 1000 | — | — |
| 2,2',6,6'-Tetrachloro[¹³ C ₁₂]biphenyl | 54L | 1000 | — | — |
| 2,3',4',5'-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 1000 | — | — |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 1000 | — | — |
| 3,4,4',5'-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 1000 | — | — |
| 2,2',3,5',6-Pentachloro[¹³ C ₁₂]biphenyl | 95L | 1000 | — | — |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 1000 | — | — |
| 2,2',4,6,6'-Pentachloro[¹³ C ₁₂]biphenyl | 104L | 1000 | — | — |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 1000 | — | — |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 1000 | — | — |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 1000 | — | — |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 1000 | — | — |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 1000 | — | — |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 1000 | — | — |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 1000 | — | — |
| 2,2',4,4',6,6'-Hexachloro[¹³ C ₁₂]biphenyl | 155L | 1000 | — | — |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 1000 | — | — |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 1000 | — | — |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 1000 | — | — |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 1000 | — | — |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 1000 | — | — |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 1000 | — | — |
| 2,2',3,4',5,6,6'-Heptachloro[¹³ C ₁₂]biphenyl | 188L | 1000 | — | — |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 1000 | — | — |
| 2,2',3,3',5,5',6,6'-Octachloro[¹³ C ₁₂]biphenyl | 202L | 1000 | — | — |
| 2,2',3,3',4,4',5,5',6-Octachloro[¹³ C ₁₂]biphenyl | 205L | 1000 | — | — |
| 2,2',3,3',4,4',5,5',6'-Nonachloro[¹³ C ₁₂]biphenyl | 208L | 1000 | — | — |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 1000 | — | — |
| RECOVERY/INTERNAL STANDARDS | | | | |
| 2,5-Dichloro[¹³ C ₁₂]biphenyl | 9L | — | 1000 | — |
| 3,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 37L | — | 1000 | — |
| 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | — | 1000 | — |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | — | 1000 | — |
| 2,3,3',4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 162L | — | 1000 | — |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | — | 1000 | — |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | — | 1000 | — |
| SAMPLING/CLEANUP STANDARDS | | | | |
| 2,3,4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 60L | — | 1000 | — |
| 2,3,3',4,5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 159L | — | 1000 | — |

PCB-CVS-A10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------------|------------------------------------|-----------------------------|
| PCB-CVS-A10-Set1 | CS1/CS3/CS5/CS7/CS9 | 1 kit (5 x 200 µl ampoules) |
| PCB-CVS-A10-Set2 | CS2/CS4/CS6/CS8/CS10 | 1 kit (5 x 200 µl ampoules) |
| PCB-CVS-A10-Set3 | CS3/CS5/CS7/CS9/CS11 | 1 kit (5 x 200 µl ampoules) |
| PCB-A10-CSL | CSL Extended Calibration/Low Level | 200 µl |
| PCB-A10-CS1 | CS1 | 200 µl |
| PCB-A10-CS2 | CS2 | 200 µl |
| PCB-A10-CS3 | CS3 | 200 µl |
| PCB-A10-CS4 | CS4 | 200 µl |

| NATIVE PCB CONGENERS | IUPAC | PCB-A10- | PCB-A10- | PCB-A10- | PCB-A10- | PCB-A10- |
|--------------------------------------|-------|----------|----------|----------|----------|----------|
| | | CSL | CS1 | CS2 | CS3 | CS4 |
| | | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) | (ng/ml) |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.05 | 0.1 | 0.2 | 0.5 | 1 |

MASS-LABELLED PCB CONGENERS

Sampling and Syringe Spikes

| | | | | | | |
|---|------|----|----|----|----|----|
| * 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 10 | 10 | 10 | 10 | 10 |
| ** 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 | 10 | 10 | 10 | 10 |
| ** 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 10 | 10 | 10 | 10 | 10 |
| ** 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 10 | 10 | 10 | 10 | 10 |

Surrogates/Extraction Spikes

| | | | | | | |
|---|------|----|----|----|----|----|
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 | 10 | 10 | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 10 | 10 | 10 | 10 | 10 |

* for pre-sampling spike

** for syringe spike / recovery standard

PCB-CVS-B10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-------------------------|---------------------------|-----------------------------|
| PCB-CVS-B10-Set1 | CS1/CS2/CS3/CS4/CS5 | 1 kit (5 x 200 µl ampoules) |
| PCB-CVS-B10-Set2 | CS2/CS3/CS4/CS5/CS6 | 1 kit (5 x 200 µl ampoules) |
| PCB-B10-CS1 | CS1 | 200 µl |
| PCB-B10-CS2 | CS2 | 200 µl |
| PCB-B10-CS3 | CS3 | 200 µl |

| NATIVE PCB CONGENERS | IUPAC | PCB-B10-CS1 (ng/ml) | PCB-B10-CS2 (ng/ml) | PCB-B10-CS3 (ng/ml) |
|---|-------|------------------------|------------------------|------------------------|
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 0.2 | 1 | 4 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 0.2 | 1 | 4 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 0.2 | 1 | 4 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 0.2 | 1 | 4 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 0.2 | 1 | 4 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 0.2 | 1 | 4 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 0.2 | 1 | 4 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 0.2 | 1 | 4 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 0.2 | 1 | 4 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 0.2 | 1 | 4 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 0.2 | 1 | 4 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 0.2 | 1 | 4 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 0.2 | 1 | 4 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 0.2 | 1 | 4 |
| MASS-LABELLED PCB CONGENERS | | | | |
| Sampling and Syringe Spikes | | | | |
| * 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 10 | 10 | 10 |
| **2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 | 10 | 10 |
| **2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 10 | 10 | 10 |
| **2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 10 | 10 | 10 |
| **2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | 10 | 10 | 10 |
| Surrogate/Extraction Spikes | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 10 | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 10 | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 10 | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 10 | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 10 | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 10 | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 10 | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 10 | 10 | 10 |

* for sampling spike

** for syringe spike

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------|----------|
| PCB-B10-CS4 | CS4 | 200 µl |
| PCB-B10-CS5 | CS5 | 200 µl |
| PCB-B10-CS6 | CS6 | 200 µl |
| PCB-B10-CS7 | CS7 | 200 µl |

| | | PCB-B10- CS4 (ng/ml) | PCB-B10- CS5 (ng/ml) | PCB-B10- CS6 (ng/ml) | PCB-B10- CS7 (ng/ml) |
|---|--------------|----------------------------|----------------------------|----------------------------|----------------------------|
| NATIVE PCB CONGENERS | | | | | |
| | IUPAC | | | | |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 20 | 100 | 400 | 1000 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 20 | 100 | 400 | 1000 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 20 | 100 | 400 | 1000 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 20 | 100 | 400 | 1000 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 20 | 100 | 400 | 1000 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 20 | 100 | 400 | 1000 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 20 | 100 | 400 | 1000 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 20 | 100 | 400 | 1000 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 20 | 100 | 400 | 1000 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 20 | 100 | 400 | 1000 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 20 | 100 | 400 | 1000 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 20 | 100 | 400 | 1000 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 20 | 100 | 400 | 1000 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 20 | 100 | 400 | 1000 |
| MASS-LABELLED PCB CONGENERS | | | | | |
| Sampling and Syringe Spikes | | | | | |
| * 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 10 | 10 | 10 | 10 |
| **2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 10 | 10 | 10 | 10 |
| **2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 10 | 10 | 10 | 10 |
| **2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 10 | 10 | 10 | 10 |
| **2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | 10 | 10 | 10 | 10 |
| Surrogate/Extraction Spikes | | | | | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 | 10 | 10 | 10 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 10 | 10 | 10 | 10 |

* for sampling spike

** for syringe spike

MASS-LABELLED PCBs: SOLUTION/MIXTURES

Support solutions for **PCB-CVS-A10**, **PCB-CVS-B10**, and **DFP-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|---|--|-------------------|---------------------|--------------------|
| PCB-LCS-A1 | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| PCB-LCS-A100 | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| PCB-LCS-A20 | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| | | PCB-LCS-A1 | PCB-LCS-A100 | PCB-LCS-A20 |
| MASS-LABELLED PCB CONGENERS | IUPAC | (ng/ml) | (ng/ml) | (ng/ml) |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 1000 | 100 | 20 |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 1000 | 100 | 20 |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 1000 | 100 | 20 |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 1000 | 100 | 20 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 1000 | 100 | 20 |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 1000 | 100 | 20 |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 1000 | 100 | 20 |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 1000 | 100 | 20 |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 1000 | 100 | 20 |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 1000 | 100 | 20 |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 1000 | 100 | 20 |
| 2,2',3,3',4,4',5-Heptachloro[¹³ C ₁₂]biphenyl | 170L | 1000 | 100 | 20 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 1000 | 100 | 20 |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 1000 | 100 | 20 |
| PCB-IS-A | Mass-Labelled PCB Solution | 1.2 ml | | |
| PCB-IS-A100 | Mass-Labelled PCB Solution | 1.2 ml | | |
| PCB-IS-A20 | Mass-Labelled PCB Solution | 1.2 ml | | |
| | | PCB-IS-A | PCB-IS-A100 | PCB-IS-A20 |
| MASS-LABELLED PCB CONGENER | IUPAC | (ng/ml) | (ng/ml) | (ng/ml) |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 1000 | 100 | 20 |
| PCB-IS-B | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| PCB-IS-B100 | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| PCB-IS-B20 | Mass-Labelled PCB Solution/Mixture | 1.2 ml | | |
| | | PCB-IS-B | PCB-IS-B100 | PCB-IS-B20 |
| MASS-LABELLED PCB CONGENERS | IUPAC | (ng/ml) | (ng/ml) | (ng/ml) |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 1000 | 100 | 20 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 1000 | 100 | 20 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 1000 | 100 | 20 |
| PCB-IS-C | Mass-Labelled PCB Syringe Spike Solution | 1.2 ml | | |
| PCB-IS-C100 | Mass-Labelled PCB Syringe Spike Solution | 1.2 ml | | |
| PCB-IS-C20 | Mass-Labelled PCB Syringe Spike Solution | 1.2 ml | | |
| | | PCB-IS-C | PCB-IS-C100 | PCB-IS-C20 |
| MASS-LABELLED PCB CONGENERS | IUPAC | (ng/ml) | (ng/ml) | (ng/ml) |
| 2,3',4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 70L | 1000 | 100 | 20 |
| 2,3,3',5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 111L | 1000 | 100 | 20 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 1000 | 100 | 20 |
| 2,2',3,3',5,5',6-Heptachloro[¹³ C ₁₂]biphenyl | 178L | 1000 | 100 | 20 |
| PCB-SS-A | Mass-Labelled PCB Solution | 1.2 ml | | |
| PCB-SS-A100 | Mass-Labelled PCB Solution | 1.2 ml | | |
| PCB-SS-A20 | Mass-Labelled PCB Solution | 1.2 ml | | |
| | | PCB-SS-A | PCB-SS-A100 | PCB-SS-A20 |
| MASS-LABELLED PCB CONGENER | IUPAC | (ng/ml) | (ng/ml) | (ng/ml) |
| 3,3',4,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 79L | 1000 | 100 | 20 |

MASS-LABELLED PCBs: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|---|----------|
| MBP-CP | Mass-Labelled Coplanar PCB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 3,3',4,4'-Tetrachloro[¹³ C ₁₂]biphenyl | 77L | 10 µg/ml |
| 3,4,4',5-Tetrachloro[¹³ C ₁₂]biphenyl | 81L | 10 µg/ml |
| 3,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 126L | 10 µg/ml |
| 3,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 169L | 10 µg/ml |
| MBP-MO | Mass-Labelled Mono-ortho PCB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 2,3,3',4,4'-Pentachloro[¹³ C ₁₂]biphenyl | 105L | 5 µg/ml |
| 2,3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 114L | 5 µg/ml |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 5 µg/ml |
| 2',3,4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 123L | 5 µg/ml |
| 2,3,3',4,4',5-Hexachloro[¹³ C ₁₂]biphenyl | 156L | 5 µg/ml |
| 2,3,3',4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 157L | 5 µg/ml |
| 2,3',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 167L | 5 µg/ml |
| 2,3,3',4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 189L | 5 µg/ml |
| MBP-CG | Mass-Labelled Mono To Decachloro PCB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 4-Chloro[¹³ C ₁₂]biphenyl | 3L | 5 µg/ml |
| 4,4'-Dichloro[¹³ C ₁₂]biphenyl | 15L | 5 µg/ml |
| 2,4',5-Trichloro[¹³ C ₁₂]biphenyl | 31L | 5 µg/ml |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 5 µg/ml |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 5 µg/ml |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 5 µg/ml |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 5 µg/ml |
| 2,2',3,3',4,4',5,5'-Octachloro[¹³ C ₁₂]biphenyl | 194L | 5 µg/ml |
| 2,2',3,3',4,4',5,5',6-Nonachloro[¹³ C ₁₂]biphenyl | 206L | 5 µg/ml |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 5 µg/ml |
| MBP-MXE | Mass-Labelled PCB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 5 µg/ml |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 5 µg/ml |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 5 µg/ml |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 5 µg/ml |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 5 µg/ml |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 5 µg/ml |
| Decachloro[¹³ C ₁₂]biphenyl | 209L | 5 µg/ml |

NATIVE PCBs: SOLUTION/MIXTURES

(*) Support solutions for **PCB-CVS-A10**, **PCB-CVS-B10**, and **DFP-CVS-B10**

| Catalogue Number | Product (nonane solution) | Qty/Conc | | |
|--------------------------------------|---------------------------|----------------------|------------------------|-----------------------|
| PCB-ST-A* | Native PCB Stock Solution | 1.2 ml | | |
| PCB-ST-A10* | Native PCB Stock Solution | 1.2 ml | | |
| PCB-ST-A2* | Native PCB Stock Solution | 1.2 ml | | |
| NATIVE PCB CONGENERS | IUPAC | PCB-ST-A* (ng/ml) | PCB-ST-A10* (ng/ml) | PCB-ST-A2* (ng/ml) |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 2000 | 10 | 2 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 2000 | 10 | 2 |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 2000 | 10 | 2 |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 2000 | 10 | 2 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 2000 | 10 | 2 |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 2000 | 10 | 2 |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 2000 | 10 | 2 |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 2000 | 10 | 2 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 2000 | 10 | 2 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 2000 | 10 | 2 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 2000 | 10 | 2 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 2000 | 10 | 2 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 2000 | 10 | 2 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 2000 | 10 | 2 |

BP-CP81

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|-----------------------------------|--------------------------------------|----------|
| BP-CP81 | Native Coplanar PCB Solution/Mixture | 1.2 ml |
| NATIVE PCB CONGENERS | IUPAC | |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 10 µg/ml |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 10 µg/ml |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 10 µg/ml |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 10 µg/ml |

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--|---|-----------|
| BP-WD | Native PCB Window Defining Solution/Mixture for DB-5 or Equivalent Column | 1.2 ml |
| NATIVE PCB CONGENERS | IUPAC | |
| Biphenyl | — | 2.5 µg/ml |
| 2-Chlorobiphenyl | 1 | 2.5 µg/ml |
| 4-Chlorobiphenyl | 3 | 2.5 µg/ml |
| 2,6-Dichlorobiphenyl | 10 | 2.5 µg/ml |
| 4,4'-Dichlorobiphenyl | 15 | 2.5 µg/ml |
| 2,2',6-Trichlorobiphenyl | 19 | 2.5 µg/ml |
| 3,4,4'-Trichlorobiphenyl | 37 | 2.5 µg/ml |
| 2,2',6,6'-Tetrachlorobiphenyl | 54 | 2.5 µg/ml |
| 3,3',4,4'-Tetrachlorobiphenyl | 77 | 2.5 µg/ml |
| 2,2',4,6,6'-Pentachlorobiphenyl | 104 | 2.5 µg/ml |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 2.5 µg/ml |
| 2,2',4,4',6,6'-Hexachlorobiphenyl | 155 | 2.5 µg/ml |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 2.5 µg/ml |
| 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 188 | 2.5 µg/ml |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 2.5 µg/ml |
| 2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 202 | 2.5 µg/ml |
| 2,3,3',4,4',5,5',6-Octachlorobiphenyl | 205 | 2.5 µg/ml |
| 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | 2.5 µg/ml |
| 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 208 | 2.5 µg/ml |
| Decachlorobiphenyl | 209 | 2.5 µg/ml |

BP-MO

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------------------------|--|----------|
| BP-MO | Solution/Mixture of Native Mono-ortho PCBs | 1.2 ml |
| NATIVE PCB CONGENERS | IUPAC | |
| 2,3,3',4,4'-Pentachlorobiphenyl | 105 | 10 µg/ml |
| 2,3,4,4',5-Pentachlorobiphenyl | 114 | 10 µg/ml |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 10 µg/ml |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 10 µg/ml |
| 2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 10 µg/ml |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 10 µg/ml |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 10 µg/ml |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 10 µg/ml |

BP-MS

BP-MS-PL1, **BP-MS-PL2**, and **BP-MS-PL3** were prepared to be used in the identification/confirmation of the PCB congeners in **BP-MS**. Although DB-5 data is provided with **BP-MS**, there can be changes in the elution order on "equivalent capillary columns".

| Catalogue Number | Product (nonane solution) | Qty/Conc | | | | |
|------------------|--|----------|--|--|--|--|
| BP-MS | Native PCB Solution/Mixture for MS Detection | 1.2 ml | | | | |
| BP-MS2 | Native PCB Solution/Mixture for MS Detection | 1.2 ml | | | | |
| BP-MS-PL1 | Native PCB Solution/Mixture for MS Detection | 1.2 ml | | | | |
| BP-MS-PL2 | Native PCB Solution/Mixture for MS Detection | 1.2 ml | | | | |
| BP-MS-PL3 | Native PCB Solution/Mixture for MS Detection | 1.2 ml | | | | |

| PCB CONGENERS | IUPAC | BP-MS (µg/ml) | BP-MS2 (µg/ml) | BP-MS- PL1 (µg/ml) | BP-MS- PL2 (µg/ml) | BP-MS- PL3 (µg/ml) |
|---|-------|------------------|-------------------|--------------------------|--------------------------|--------------------------|
| 2-Chlorobiphenyl | 1 | 2.0 | — | — | — | — |
| 4-Chlorobiphenyl | 3 | 2.0 | — | — | — | — |
| 2,2'-Dichlorobiphenyl | 4 | 2.0 | — | — | — | — |
| 2,4'-Dichlorobiphenyl | 8 | 2.0 | — | — | — | — |
| 2,6-Dichlorobiphenyl | 10 | 2.0 | — | — | — | — |
| *4,4'-Dichlorobiphenyl | 15 | 2.0 | — | — | — | — |
| *2,2',5-Trichlorobiphenyl | 18 | 2.0 | — | — | — | — |
| 2,2',6-Trichlorobiphenyl | 19 | 2.0 | — | — | — | — |
| 2,3,4'-Trichlorobiphenyl | 22 | 2.0 | — | — | — | — |
| 2,4,4'-Trichlorobiphenyl | 28 | 2.0 | — | — | — | — |
| 2',3,4-Trichlorobiphenyl | 33 | 2.0 | — | — | — | — |
| 3,4,4'-Trichlorobiphenyl | 37 | 2.0 | — | — | — | — |
| 2,2',3,3'-Tetrachlorobiphenyl | 40 | — | 2.0 | — | — | — |
| 2,2',3,4-Tetrachlorobiphenyl | 41 | — | 2.0 | — | — | — |
| *2,2',3,5'-Tetrachlorobiphenyl | 44 | 2.0 | — | — | — | — |
| *2,2',4,5'-Tetrachlorobiphenyl | 49 | 2.0 | — | — | 2.0 | — |
| *2,2',5,5'-Tetrachlorobiphenyl | 52 | 2.0 | — | 2.0 | — | — |
| *2,2',6,6'-Tetrachlorobiphenyl | 54 | 2.0 | — | — | — | — |
| 2,3,4,4'-Tetrachlorobiphenyl | 60 | — | 2.0 | — | — | — |
| 2,3',4,4'-Tetrachlorobiphenyl | 66 | — | 2.0 | — | — | — |
| 2,3',4',5-Tetrachlorobiphenyl | 70 | 2.0 | — | 2.0 | — | — |
| 2,4,4',5-Tetrachlorobiphenyl | 74 | 2.0 | — | — | 2.0 | — |
| *3,3',4,4'-Tetrachlorobiphenyl | 77 | 2.0 | — | — | — | 2.0 |
| 3,4,4',5-Tetrachlorobiphenyl | 81 | 2.0 | — | — | — | — |
| *2,2',3,4,5'-Pentachlorobiphenyl | 87 | 2.0 | — | 2.0 | — | — |
| 2,2',3,4',5-Pentachlorobiphenyl | 90 | — | 2.0 | — | — | — |
| 2,2',3,5',6-Pentachlorobiphenyl | 95 | 2.0 | — | — | 2.0 | — |
| 2,2',4,4',5-Pentachlorobiphenyl | 99 | 2.0 | — | — | 2.0 | — |
| *2,2',4,5,5'-Pentachlorobiphenyl | 101 | 2.0 | — | 2.0 | — | — |
| 2,2',4,6,6'-Pentachlorobiphenyl | 104 | 2.0 | — | — | — | — |
| *2,3,3',4,4'-Pentachlorobiphenyl | 105 | 2.0 | — | — | — | — |
| 2,3,3',4',6-Pentachlorobiphenyl | 110 | 2.0 | — | 2.0 | — | — |
| *2,3,4,4',5-Pentachlorobiphenyl | 114 | 2.0 | — | — | — | — |
| *2,3',4,4',5-Pentachlorobiphenyl | 118 | 2.0 | — | — | — | — |
| 2,3',4,4',6-Pentachlorobiphenyl | 119 | 2.0 | — | — | — | — |
| 2',3,4,4',5-Pentachlorobiphenyl | 123 | 2.0 | — | — | — | — |
| 3,3',4,4',5-Pentachlorobiphenyl | 126 | 2.0 | — | — | — | — |
| *2,2',3,3',4,4'-Hexachlorobiphenyl | 128 | 2.0 | — | — | — | 2.0 |
| 2,2',3,3',4,5-Hexachlorobiphenyl | 129 | — | 2.0 | — | — | — |
| 2,2',3,4,4',5-Hexachlorobiphenyl | 137 | — | 2.0 | — | — | — |
| *2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | 2.0 | — | 2.0 | — | — |
| 2,2',3,4,5,5'-Hexachlorobiphenyl | 141 | — | 2.0 | — | — | — |
| 2,2',3,4',5,6-Hexachlorobiphenyl | 149 | 2.0 | — | — | 2.0 | — |
| *2,2',3,5,5',6-Hexachlorobiphenyl | 151 | 2.0 | — | — | 2.0 | — |
| *2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | 2.0 | — | 2.0 | — | — |
| 2,2',4,4',6,6'-Hexachlorobiphenyl | 155 | 2.0 | — | 2.0 | — | — |
| *2,3,3',4,4',5-Hexachlorobiphenyl | 156 | 2.0 | — | — | — | — |
| 2,3,3',4,4',5'-Hexachlorobiphenyl | 157 | 2.0 | — | — | — | — |
| 2,3,3',4,4',6-Hexachlorobiphenyl | 158 | 2.0 | — | — | 2.0 | — |
| 2,3',4,4',5,5'-Hexachlorobiphenyl | 167 | 2.0 | — | — | — | — |
| 2,3',4,4',5',6-Hexachlorobiphenyl | 168 | 2.0 | — | — | 2.0 | — |
| 3,3',4,4',5,5'-Hexachlorobiphenyl | 169 | 2.0 | — | — | — | — |
| *2,2',3,3',4,4',5-Heptachlorobiphenyl | 170 | 2.0 | — | — | — | — |
| *2,2',3,3',4,4',6-Heptachlorobiphenyl | 171 | 2.0 | — | — | 2.0 | — |
| 2,2',3,3',4',5,6-Heptachlorobiphenyl | 177 | 2.0 | — | 2.0 | — | — |
| 2,2',3,3',5,5',6-Heptachlorobiphenyl | 178 | 2.0 | — | — | — | 2.0 |
| *2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 2.0 | — | 2.0 | — | — |
| *2,2',3,4,4',5',6-Heptachlorobiphenyl | 183 | 2.0 | — | — | — | — |
| *2,2',3,4',5,5',6-Heptachlorobiphenyl | 187 | 2.0 | — | — | — | — |
| 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 188 | 2.0 | — | 2.0 | — | — |
| *2,3,3',4,4',5,5'-Heptachlorobiphenyl | 189 | 2.0 | — | — | — | — |
| *2,3,3',4,4',5',6-Heptachlorobiphenyl | 191 | 2.0 | — | — | — | — |
| 2,3,3',4',5,5',6-Heptachlorobiphenyl | 193 | — | 2.0 | — | — | — |
| *2,2',3,3',4,4',5,5'-Octachlorobiphenyl | 194 | 2.0 | — | — | — | — |
| *2,2',3,3',4,5,5',6'-Octachlorobiphenyl | 199 | 2.0 | — | — | — | — |
| *2,2',3,3',4,5',6,6'-Octachlorobiphenyl | 201 | 2.0 | — | 2.0 | — | — |
| *2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 202 | 2.0 | — | — | — | 2.0 |
| 2,2',3,4,4',5,5',6-Octachlorobiphenyl | 203 | — | 2.0 | — | — | — |
| *2,3,3',4,4',5,5',6-Octachlorobiphenyl | 205 | 2.0 | — | — | — | — |
| *2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 206 | 2.0 | — | — | — | — |
| *2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 208 | 2.0 | — | — | — | — |
| *Decachlorobiphenyl | 209 | 2.0 | — | — | — | — |

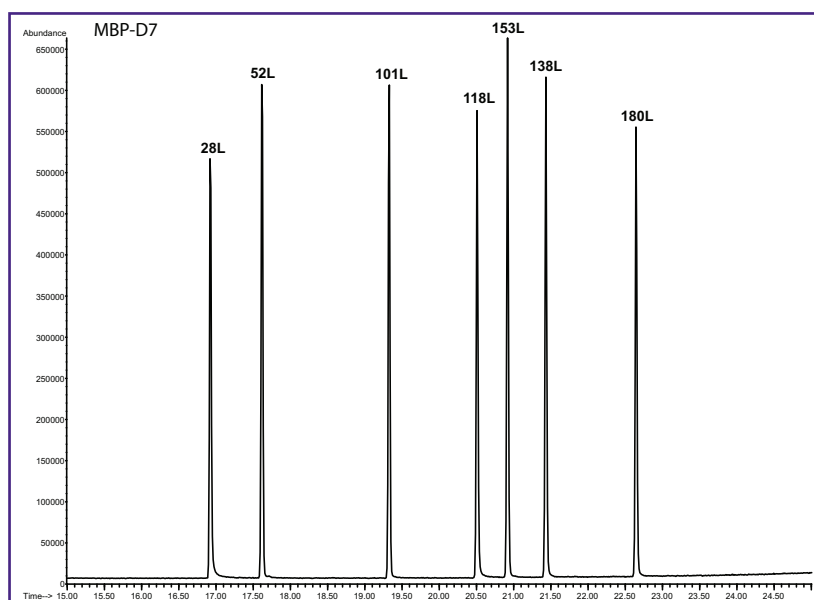
Congeners marked with an asterisk (*) are concentration-certified by direct comparison to the NRCC CLB-1 solutions.

Solution/Mixtures for the analysis of the Dutch 7 PCB Congeners.

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| BP-D7 | Native PCB Congener Solution/Mixture | 1.2 ml |
| MBP-D7 | Mass-Labelled PCB Congener Solution/Mixture | 1.2 ml |

| NATIVE PCB CONGENERS | IUPAC | BP-D7 (µg/ml) |
|--------------------------------------|-------|---------------|
| 2,4,4'-Trichlorobiphenyl | 28 | 10 |
| 2,2',5,5'-Tetrachlorobiphenyl | 52 | 10 |
| 2,2',4,5,5'-Pentachlorobiphenyl | 101 | 10 |
| 2,3',4,4',5-Pentachlorobiphenyl | 118 | 10 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl | 138 | 10 |
| 2,2',4,4',5,5'-Hexachlorobiphenyl | 153 | 10 |
| 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 180 | 10 |

| MASS-LABELLED PCB CONGENERS | IUPAC | MBP-D7 (µg/ml) |
|---|-------|----------------|
| 2,4,4'-Trichloro[¹³ C ₁₂]biphenyl | 28L | 5.0 |
| 2,2',5,5'-Tetrachloro[¹³ C ₁₂]biphenyl | 52L | 5.0 |
| 2,2',4,5,5'-Pentachloro[¹³ C ₁₂]biphenyl | 101L | 5.0 |
| 2,3',4,4',5-Pentachloro[¹³ C ₁₂]biphenyl | 118L | 5.0 |
| 2,2',3,4,4',5'-Hexachloro[¹³ C ₁₂]biphenyl | 138L | 5.0 |
| 2,2',4,4',5,5'-Hexachloro[¹³ C ₁₂]biphenyl | 153L | 5.0 |
| 2,2',3,4,4',5,5'-Heptachloro[¹³ C ₁₂]biphenyl | 180L | 5.0 |



HRGC/LRMS Data: MBP-D7 on a 30m DB-5 column.

MASS-LABELLED PCDDs/PCDFs/PCBs: SOLUTION/MIXTURES

These three solutions were designed and prepared as support solutions to be used with the following calibration sets:

DF-CVS-A10 (see Page 42)

DF-CVS-B10 (see Page 44)

as well as:

PCB-CVS-A10

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|--|----------|
| DFP-LCS-A | Mass-Labelled PCDD/PCDF/PCB Solution/Mixture | 1.2 ml |
| MASS-LABELLED PCDDs | | |
| 2,3,7,8-Tetrachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,7,8-Pentachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,4,7,8-Hexachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,6,7,8-Hexachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,7,8,9-Hexachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| 1,2,3,4,6,7,8-Heptachloro ^[13C] dibenzo-p-dioxin | | 10 ng/ml |
| Octachloro ^[13C] dibenzo-p-dioxin | | 20 ng/ml |
| MASS-LABELLED PCDFs | | |
| 2,3,7,8-Tetrachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,7,8-Pentachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 2,3,4,7,8-Pentachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,4,7,8-Hexachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,6,7,8-Hexachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,7,8,9-Hexachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 2,3,4,6,7,8-Hexachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,4,6,7,8-Heptachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,4,7,8,9-Heptachloro ^[13C] dibenzofuran | | 10 ng/ml |
| Octachloro ^[13C] dibenzofuran | | 20 ng/ml |
| MASS-LABELLED PCBs | | |
| | IUPAC | |
| 3,3',4,4'-Tetrachloro ^[13C] biphenyl | 77L | 10 ng/ml |
| 3,4,4',5-Tetrachloro ^[13C] biphenyl | 81L | 10 ng/ml |
| 2,3,3',4,4'-Pentachloro ^[13C] biphenyl | 105L | 10 ng/ml |
| 2,3,4,4',5-Pentachloro ^[13C] biphenyl | 114L | 10 ng/ml |
| 2,3',4,4',5-Pentachloro ^[13C] biphenyl | 118L | 10 ng/ml |
| 2',3,4,4',5-Pentachloro ^[13C] biphenyl | 123L | 10 ng/ml |
| 3,3',4,4',5-Pentachloro ^[13C] biphenyl | 126L | 10 ng/ml |
| 2,3,3',4,4',5-Hexachloro ^[13C] biphenyl | 156L | 10 ng/ml |
| 2,3,3',4,4',5'-Hexachloro ^[13C] biphenyl | 157L | 10 ng/ml |
| 2,3',4,4',5,5'-Hexachloro ^[13C] biphenyl | 167L | 10 ng/ml |
| 3,3',4,4',5,5'-Hexachloro ^[13C] biphenyl | 169L | 10 ng/ml |
| 2,2',3,3',4,4',5-Heptachloro ^[13C] biphenyl | 170L | 10 ng/ml |
| 2,2',3,4,4',5,5'-Heptachloro ^[13C] biphenyl | 180L | 10 ng/ml |
| 2,3,3',4,4',5,5'-Heptachloro ^[13C] biphenyl | 189L | 10 ng/ml |
| DFP-IS-A | Mass-Labelled PCDF/PCB Syringe Spike | 1.2 ml |
| | IUPAC | |
| 2,3',4,5-Tetrachloro ^[13C] biphenyl | 70L | 10 ng/ml |
| 1,2,3,4,6,9-Hexachloro ^[13C] dibenzofuran | | 10 ng/ml |
| 1,2,3,4,6,8,9-Heptachloro ^[13C] dibenzofuran | | 10 ng/ml |
| DFP-SS-A | Mass-Labelled PCDD/PCB Sampling Spike | 1.2 ml |
| | IUPAC | |
| 3,3',4,5'-Tetrachloro ^[13C] biphenyl | 79L | 50 ng/ml |
| 1,2,3,4-Tetrachloro ^[13C] dibenzo-p-dioxin | | 50 ng/ml |

PBDEs & PBBs: POLYBROMINATED DIPHENYL ETHERS (PBDES) & POLYBROMINATED BIPHENYLS (PBBs)

This section is devoted to polybrominated diphenyl ethers (PBDEs) and polybrominated biphenyls (PBBs) and contains the PBDE calibration solutions listed below:

BFR-CVS calibration set and support solutions
BDE-CVS-F calibration set and support solutions
BDE-CVS-G calibration set and support solutions

Also included in this section are individual native and mass-labelled PBDEs and PBBs as well as selected technical mixtures and individual PBDE metabolites:

Native and mass-labelled PBDEs
Mass-labelled hydroxy-PBDEs
Native and mass-labelled methoxy-PBDEs
Native and mass-labelled PBBs



BFR-CVS

| Catalogue Number | Product (toluene solution) | Qty/Conc | | | | |
|--|--|--------------------|--------------------|--------------------|--------------------|--------------------|
| BFR-CVS | Polybrominated Diphenyl Ethers/Brominated Flame Retardants | 1 kit | | | | |
| | Calibration Solutions CS1-CS5 | (5 ampoules) | | | | |
| BFR-CS1, BFR-CS2, BFR-CS3, BFR-CS4, BFR-CS5 | Individual Calibration Solutions | 200 µl each | | | | |
| | | BFR-CS1 (ng/ml) | BFR-CS2 (ng/ml) | BFR-CS3 (ng/ml) | BFR-CS4 (ng/ml) | BFR-CS5 (ng/ml) |
| NATIVE PBDEs/BFRs | | | | | | |
| 2-Bromodiphenyl ether | BDE-1 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 3-Bromodiphenyl ether | BDE-2 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 4-Bromodiphenyl ether | BDE-3 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 2,4-Dibromodiphenyl ether | BDE-7 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 2,6-Dibromodiphenyl ether | BDE-10 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 4,4'-Dibromodiphenyl ether | BDE-15 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 2,2',4-Tribromodiphenyl ether | BDE-17 (>96%) | 0.24 | 0.96 | 4.8 | 19 | 96 |
| 2,4,4'-Tribromodiphenyl ether | BDE-28 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 2,4,6-Tribromodiphenyl ether | BDE-30 | 0.25 | 1.0 | 5.0 | 20 | 100 |
| Pentabromoethylbenzene | PBEB | 0.25 | 1.0 | 5.0 | 20 | 100 |
| Hexabromobenzene | HBBZ | 0.25 | 1.0 | 5.0 | 20 | 100 |
| 2,2',4,4'-Tetrabromodiphenyl ether | BDE-47 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,5'-Tetrabromodiphenyl ether | BDE-49 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3',4,4'-Tetrabromodiphenyl ether | BDE-66 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3',4',6-Tetrabromodiphenyl ether | BDE-71 | 0.5 | 2.0 | 10 | 40 | 200 |
| 3,3',4,4'-Tetrabromodiphenyl ether | BDE-77 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | BDE-85 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,4',5-Pentabromodiphenyl ether | BDE-99 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,4',6-Pentabromodiphenyl ether | BDE-100 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3',4,4',6-Pentabromodiphenyl ether | BDE-119 | 0.5 | 2.0 | 10 | 40 | 200 |
| 3,3',4,4',5-Pentabromodiphenyl ether | BDE-126 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether | BDE-138 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',3,4,4',6-Hexabromodiphenyl ether | BDE-139 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',3,4,4',6'-Hexabromodiphenyl ether | BDE-140 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | BDE-153 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | BDE-154 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3,3',4,4',5-Hexabromodiphenyl ether | BDE-156 | 0.5 | 2.0 | 10 | 40 | 200 |
| 3,3',4,4',5,5'-Hexabromodiphenyl ether | BDE-169 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',4,4',5,5'-Hexabromobiphenyl | BB-153 | 0.5 | 2.0 | 10 | 40 | 200 |
| 1,2-Bis(2,4,6-tribromophenoxy)ethane | BTBPE | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,2',3,3',4,4',6-Heptabromodiphenyl ether | BDE-171 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,4,4',5,5'-Heptabromodiphenyl ether | BDE-180 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | BDE-183 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,4,4',6,6'-Heptabromodiphenyl ether | BDE-184 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,3,3',4,4',5',6-Heptabromodiphenyl ether | BDE-191 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | BDE-196 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | BDE-197 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,3',4,5',6,6'-Octabromodiphenyl ether | BDE-201 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,4,4',5,5',6-Octabromodiphenyl ether | BDE-203 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether | BDE-204 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,3,3',4,4',5,5',6-Octabromodiphenyl ether | BDE-205 | 1.0 | 4.0 | 20 | 80 | 400 |
| 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | BDE-206 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | BDE-207 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,2',3,3',4,5,5',6,6'-Nonabromodiphenyl ether | BDE-208 | 2.5 | 10 | 50 | 200 | 1000 |
| Decabromodiphenyl ether | BDE-209 | 2.5 | 10 | 50 | 200 | 1000 |
| Decabromodiphenylethane | DBDPE | 5.0 | 20 | 100 | 400 | 2000 |
| BFR-LCS | | | | | | |
| 4-Bromo[¹³ C ₁₂]diphenyl ether | MBDE-3 | 25 | 25 | 25 | 25 | 25 |
| 4,4'-Dibromo[¹³ C ₁₂]diphenyl ether | MBDE-15 | 25 | 25 | 25 | 25 | 25 |
| 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | MBDE-28 | 25 | 25 | 25 | 25 | 25 |
| Hexabromo[¹³ C ₁₂]benzene | MHBBZ | 25 | 25 | 25 | 25 | 25 |
| 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | MBDE-47 | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | MBDE-77 | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | MBDE-99 | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',6-Pentabromo[¹³ C ₁₂]diphenyl ether | MBDE-100 | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | MBDE-126 | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | MBDE-153 | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | MBDE-154 | 50 | 50 | 50 | 50 | 50 |
| 3,3',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | MBDE-169 | 50 | 50 | 50 | 50 | 50 |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]biphenyl | MBB-153 | 50 | 50 | 50 | 50 | 50 |
| 1,2-Bis(2,4,6-tribromo[¹³ C ₁₂]phenoxy)ethane | MBTBPE | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',5',6-Heptabromo[¹³ C ₁₂]diphenyl ether | MBDE-183 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',6,6'-Octabromo[¹³ C ₁₂]diphenyl ether | MBDE-197 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5',6-Octabromo[¹³ C ₁₂]diphenyl ether | MBDE-205 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,6,6'-Nonabromo[¹³ C ₁₂]diphenyl ether | MBDE-207 | 250 | 250 | 250 | 250 | 250 |
| Decabromo[¹³ C ₁₂]diphenyl ether | MBDE-209 | 250 | 250 | 250 | 250 | 250 |
| Decabromo[¹³ C ₁₂]diphenylethane | MDBDPE | 500 | 500 | 500 | 500 | 500 |
| BFR-ISS | | | | | | |
| 3,3',4,5'-Tetrabromo[¹³ C ₁₂]diphenyl ether | MBDE-79 | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',6-Hexabromo[¹³ C ₁₂]diphenyl ether | MBDE-139 | 50 | 50 | 50 | 50 | 50 |
| 2,2',3,4,4',5,5'-Heptabromo[¹³ C ₁₂]diphenyl ether | MBDE-180 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5',6-Nonabromo[¹³ C ₁₂]diphenyl ether | MBDE-206 | 250 | 250 | 250 | 250 | 250 |
| BFR-SCS | | | | | | |
| 2,2',3,4,4',5'-Hexabromo[¹³ C ₁₂]diphenyl ether | MBDE-138 | 50 | 50 | 50 | 50 | 50 |

| Catalogue Number | Product (toluene solution) | Qty/Conc | | | |
|--|---|--------------------|--------------------|--------------------|--------------------|
| BFR-LCS | Labelled Compounds Stock Solution | 1.2 ml | | | |
| BFR-ISS | Internal/Injection Standards | 1.2 ml | | | |
| BFR-SCS | Sampling/Cleanup Standard | 1.2 ml | | | |
| BFR-PAR | Native Compounds Stock Solution (nonane/toluene solution) | 1.2 ml | | | |
| | | BFR-LCS (ng/ml) | BFR-ISS (ng/ml) | BFR-SCS (ng/ml) | BFR-PAR (ng/ml) |
| NATIVE PBDEs/BFRs | | | | | |
| 2-Bromodiphenyl ether | BDE-1 | — | — | — | 200 |
| 3-Bromodiphenyl ether | BDE-2 | — | — | — | 200 |
| 4-Bromodiphenyl ether | BDE-3 | — | — | — | 200 |
| 2,4-Dibromodiphenyl ether | BDE-7 | — | — | — | 200 |
| 2,6-Dibromodiphenyl ether | BDE-10 | — | — | — | 200 |
| 4,4'-Dibromodiphenyl ether | BDE-15 | — | — | — | 200 |
| 2,2',4-Tribromodiphenyl ether | BDE-17 (>96%) | — | — | — | 192 |
| 2,4,4'-Tribromodiphenyl ether | BDE-28 | — | — | — | 200 |
| 2,4,6-Tribromodiphenyl ether | BDE-30 | — | — | — | 200 |
| Pentabromoethylbenzene | PBEB | — | — | — | 200 |
| Hexabromobenzene | HBBZ | — | — | — | 200 |
| 2,2',4,4'-Tetrabromodiphenyl ether | BDE-47 | — | — | — | 400 |
| 2,2',4,5'-Tetrabromodiphenyl ether | BDE-49 | — | — | — | 400 |
| 2,3',4,4'-Tetrabromodiphenyl ether | BDE-66 | — | — | — | 400 |
| 2,3',4',6-Tetrabromodiphenyl ether | BDE-71 | — | — | — | 400 |
| 3,3',4,4'-Tetrabromodiphenyl ether | BDE-77 | — | — | — | 400 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | BDE-85 | — | — | — | 400 |
| 2,2',4,4',5-Pentabromodiphenyl ether | BDE-99 | — | — | — | 400 |
| 2,2',4,4',6-Pentabromodiphenyl ether | BDE-100 | — | — | — | 400 |
| 2,3',4,4',6-Pentabromodiphenyl ether | BDE-119 | — | — | — | 400 |
| 3,3',4,4',5-Pentabromodiphenyl ether | BDE-126 | — | — | — | 400 |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether | BDE-138 | — | — | — | 400 |
| 2,2',3,4,4',6-Hexabromodiphenyl ether | BDE-139 | — | — | — | 400 |
| 2,2',3,4,4',6'-Hexabromodiphenyl ether | BDE-140 | — | — | — | 400 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | BDE-153 | — | — | — | 400 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | BDE-154 | — | — | — | 400 |
| 2,3,3',4,4',5-Hexabromodiphenyl ether | BDE-156 | — | — | — | 400 |
| 3,3',4,4',5,5'-Hexabromodiphenyl ether | BDE-169 | — | — | — | 400 |
| 2,2',4,4',5,5'-Hexabromobiphenyl | BB-153 | — | — | — | 400 |
| 1,2-Bis(2,4,6-tribromophenoxy)ethane | BTBPE | — | — | — | 400 |
| 2,2',3,3',4,4',6-Heptabromodiphenyl ether | BDE-171 | — | — | — | 800 |
| 2,2',3,4,4',5,5'-Heptabromodiphenyl ether | BDE-180 | — | — | — | 800 |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | BDE-183 | — | — | — | 800 |
| 2,2',3,4,4',6,6'-Heptabromodiphenyl ether | BDE-184 | — | — | — | 800 |
| 2,3,3',4,4',5',6-Heptabromodiphenyl ether | BDE-191 | — | — | — | 800 |
| 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | BDE-196 | — | — | — | 800 |
| 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | BDE-197 | — | — | — | 800 |
| 2,2',3,3',4,5',6,6'-Octabromodiphenyl ether | BDE-201 | — | — | — | 800 |
| 2,2',3,4,4',5,5',6-Octabromodiphenyl ether | BDE-203 | — | — | — | 800 |
| 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether | BDE-204 | — | — | — | 800 |
| 2,3,3',4,4',5,5',6-Octabromodiphenyl ether | BDE-205 | — | — | — | 800 |
| 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | BDE-206 | — | — | — | 2000 |
| 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | BDE-207 | — | — | — | 2000 |
| 2,2',3,3',4,5,5',6,6'-Nonabromodiphenyl ether | BDE-208 | — | — | — | 2000 |
| Decabromodiphenyl ether | BDE-209 | — | — | — | 2000 |
| Decabromodiphenylethane | DBDPE | — | — | — | 4000 |
| BFR-LCS | | | | | |
| 4-Bromo[¹³ C] ₁₂ diphenyl ether | MBDE-3 | 100 | — | — | — |
| 4,4'-Dibromo[¹³ C] ₁₂ diphenyl ether | MBDE-15 | 100 | — | — | — |
| 2,4,4'-Tribromo[¹³ C] ₁₂ diphenyl ether | MBDE-28 | 100 | — | — | — |
| Hexabromo[¹³ C] ₁₂ benzene | MHBBZ | 100 | — | — | — |
| 2,2',4,4'-Tetrabromo[¹³ C] ₁₂ diphenyl ether | MBDE-47 | 200 | — | — | — |
| 3,3',4,4'-Tetrabromo[¹³ C] ₁₂ diphenyl ether | MBDE-77 | 200 | — | — | — |
| 2,2',4,4',5-Pentabromo[¹³ C] ₁₂ diphenyl ether | MBDE-99 | 200 | — | — | — |
| 2,2',4,4',6-Pentabromo[¹³ C] ₁₂ diphenyl ether | MBDE-100 | 200 | — | — | — |
| 3,3',4,4',5-Pentabromo[¹³ C] ₁₂ diphenyl ether | MBDE-126 | 200 | — | — | — |
| 2,2',4,4',5,5'-Hexabromo[¹³ C] ₁₂ diphenyl ether | MBDE-153 | 200 | — | — | — |
| 2,2',4,4',5,6'-Hexabromo[¹³ C] ₁₂ diphenyl ether | MBDE-154 | 200 | — | — | — |
| 3,3',4,4',5,5'-Hexabromo[¹³ C] ₁₂ diphenyl ether | MBDE-169 | 200 | — | — | — |
| 2,2',4,4',5,5'-Hexabromo[¹³ C] ₁₂ biphenyl | MBB-153 | 200 | — | — | — |
| 1,2-Bis(2,4,6-tribromo[¹³ C] ₁₂ phenoxy)ethane | MBTBPPE | 200 | — | — | — |
| 2,2',3,4,4',5',6-Heptabromo[¹³ C] ₁₂ diphenyl ether | MBDE-183 | 400 | — | — | — |
| 2,2',3,3',4,4',6,6'-Octabromo[¹³ C] ₁₂ diphenyl ether | MBDE-197 | 400 | — | — | — |
| 2,3,3',4,4',5,5',6-Octabromo[¹³ C] ₁₂ diphenyl ether | MBDE-205 | 400 | — | — | — |
| 2,2',3,3',4,4',5,6,6'-Nonabromo[¹³ C] ₁₂ diphenyl ether | MBDE-207 | 1000 | — | — | — |
| Decabromo[¹³ C] ₁₂ diphenyl ether | MBDE-209 | 1000 | — | — | — |
| Decabromo[¹³ C] ₁₂ diphenylethane | MDBDPE | 2000 | — | — | — |
| BFR-ISS | | | | | |
| 3,3',4,5'-Tetrabromo[¹³ C] ₁₂ diphenyl ether | MBDE-79 | — | 200 | — | — |
| 2,2',3,4,4',6-Hexabromo[¹³ C] ₁₂ diphenyl ether | MBDE-139 | — | 200 | — | — |
| 2,2',3,4,4',5,5'-Heptabromo[¹³ C] ₁₂ diphenyl ether | MBDE-180 | — | 400 | — | — |
| 2,2',3,3',4,4',5,5',6-Nonabromo[¹³ C] ₁₂ diphenyl ether | MBDE-206 | — | 1000 | — | — |
| BFR-SCS | | | | | |
| 2,2',3,4,4',5'-Hexabromo[¹³ C] ₁₂ diphenyl ether | MBDE-138 | — | — | 400 | — |

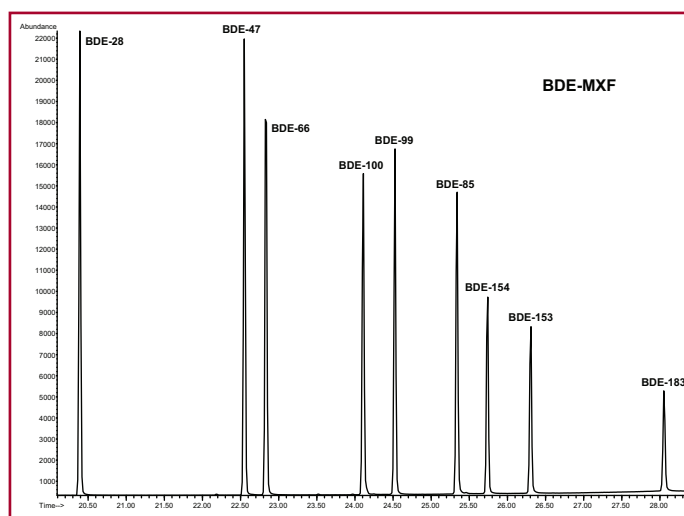
BDE-CVS-F

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|-----------------------|
| BDE-CVS-F | BDE-CVS-F Calibration Solutions CS1-CS5 | 1 kit (5 ampoules) |
| BDE-CS1-F | CS1 | 200 µl |
| BDE-CS2-F | CS2 | 200 µl |
| BDE-CS3-F | CS3 | 200 µl |
| BDE-CS4-F | CS4 | 200 µl |
| BDE-CS5-F | CS5 | 200 µl |

| NATIVE PBDE CONGENERS | IUPAC | BDE- CS1-F (ng/ml) | BDE- CS2-F (ng/ml) | BDE- CS3-F (ng/ml) | BDE- CS4-F (ng/ml) | BDE- CS5-F (ng/ml) |
|---|-------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| 2,4,4'-Tribromodiphenyl ether | 28 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',4,4'-Tetrabromodiphenyl ether | 47 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,3',4,4'-Tetrabromodiphenyl ether | 66 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | 85 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',4,4',5-Pentabromodiphenyl ether | 99 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',4,4',6-Pentabromodiphenyl ether | 100 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | 153 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | 154 | 1.0 | 5.0 | 25 | 100 | 500 |
| 2,2',3,4,4',5,6'-Heptabromodiphenyl ether | 183 | 1.0 | 5.0 | 25 | 100 | 500 |

| MASS-LABELLED PBDE SURROGATE STANDARDS | IUPAC | BDE- CS1-F (ng/ml) | BDE- CS2-F (ng/ml) | BDE- CS3-F (ng/ml) | BDE- CS4-F (ng/ml) | BDE- CS5-F (ng/ml) |
|--|-------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 28L | 20 | 20 | 20 | 20 | 20 |
| 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 47L | 20 | 20 | 20 | 20 | 20 |
| 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 99L | 20 | 20 | 20 | 20 | 20 |
| 2,2',4,4',6-Pentabromo[¹³ C ₁₂]diphenyl ether | 100L | 20 | 20 | 20 | 20 | 20 |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 153L | 20 | 20 | 20 | 20 | 20 |
| 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 154L | 20 | 20 | 20 | 20 | 20 |
| 2,2',3,4,4',5,6'-Heptabromo[¹³ C ₁₂]diphenyl ether | 183L | 20 | 20 | 20 | 20 | 20 |

| MASS-LABELLED PBDE RECOVERY STANDARDS | IUPAC | BDE- CS1-F (ng/ml) | BDE- CS2-F (ng/ml) | BDE- CS3-F (ng/ml) | BDE- CS4-F (ng/ml) | BDE- CS5-F (ng/ml) |
|---|-------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| 3,3',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 77L | 20 | 20 | 20 | 20 | 20 |
| 2,2',3,4,4',5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 138L | 20 | 20 | 20 | 20 | 20 |



BDE-MXF; HRGC/LRMS TIC Chromatogram

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|------------------------------------|----------|
| MBDE-MXFS | Mass-Labelled PBDE Surrogate Stock | 1.2 ml |
| MBDE-MXFR | Mass-Labelled PBDE Recovery Stock | 1.2 ml |
| BDE-MXF | Native PBDE Stock | 1.2 ml |

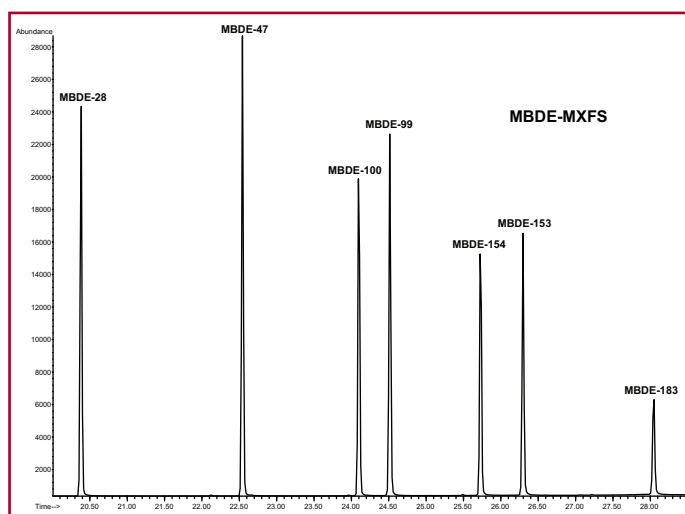
| NATIVE PBDE CONGENERS | IUPAC | MBDE-MXFS (ng/ml) | MBDE-MXFR (ng/ml) | BDE-MXF (ng/ml) |
|---|-------|-------------------|-------------------|-----------------|
| 2,4,4'-Tribromodiphenyl ether | 28 | — | — | 2000 |
| 2,2',4,4'-Tetrabromodiphenyl ether | 47 | — | — | 2000 |
| 2,3',4,4'-Tetrabromodiphenyl ether | 66 | — | — | 2000 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | 85 | — | — | 2000 |
| 2,2',4,4',5-Pentabromodiphenyl ether | 99 | — | — | 2000 |
| 2,2',4,4',6-Pentabromodiphenyl ether | 100 | — | — | 2000 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | 153 | — | — | 2000 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | 154 | — | — | 2000 |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | 183 | — | — | 2000 |

MASS-LABELLED PBDE SURROGATE STANDARDS

| | | | | |
|--|------|------|---|---|
| 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 28L | 2000 | — | — |
| 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 47L | 2000 | — | — |
| 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 99L | 2000 | — | — |
| 2,2',4,4',6-Pentabromo[¹³ C ₁₂]diphenyl ether | 100L | 2000 | — | — |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 153L | 2000 | — | — |
| 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 154L | 2000 | — | — |
| 2,2',3,4,4',5',6-Heptabromo[¹³ C ₁₂]diphenyl ether | 183L | 2000 | — | — |

MASS-LABELLED PBDE RECOVERY STANDARDS

| | | | | |
|---|------|---|------|---|
| 3,3',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 77L | — | 2000 | — |
| 2,2',3,4,4',5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 138L | — | 2000 | — |



MBDE-MXFS; HRGC/LRMS TIC Chromatogram

BDE-CVS-G

| Catalogue Number | Product (nonane/toluene solution) | Qty/Conc |
|------------------|--|-----------------------|
| BDE-CVS-G | BDE-CVS-G Calibration Solutions CS1-CS5 | 1 kit (5 ampoules) |
| BDE-CS1-G | CS1 | 200 µl |
| BDE-CS2-G | CS2 | 200 µl |
| BDE-CS3-G | CS3 | 200 µl |
| BDE-CS4-G | CS4 | 200 µl |
| BDE-CS5-G | CS5 | 200 µl |

NOTE: This set of calibration solutions were designed to be used with BDE-MXE as the native PBDE stock solution.

| NATIVE PBDE CONGENERS | IUPAC | BDE- CS1-G (ng/ml) | BDE- CS2-G (ng/ml) | BDE- CS3-G (ng/ml) | BDE- CS4-G (ng/ml) | BDE- CS5-G (ng/ml) |
|--|-------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| 4-Bromodiphenyl ether | 3 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,4-Dibromodiphenyl ether | 7 | 1.0 | 5.0 | 20 | 100 | 400 |
| 4,4'-Dibromodiphenyl ether | 15 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',4-Tribromodiphenyl ether (>96%) | 17 | 0.96 | 4.8 | 19.2 | 96 | 384 |
| 2,4,4'-Tribromodiphenyl ether | 28 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',4,4'-Tetrabromodiphenyl ether | 47 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',4,5'-Tetrabromodiphenyl ether | 49 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,3',4,4'-Tetrabromodiphenyl ether | 66 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,3',4',6-Tetrabromodiphenyl ether | 71 | 1.0 | 5.0 | 20 | 100 | 400 |
| 3,3',4,4'-Tetrabromodiphenyl ether | 77 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | 85 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',4,4',5-Pentabromodiphenyl ether | 99 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',4,4',6-Pentabromodiphenyl ether | 100 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,3',4,4',6-Pentabromodiphenyl ether | 119 | 1.0 | 5.0 | 20 | 100 | 400 |
| 3,3',4,4',5-Pentabromodiphenyl ether | 126 | 1.0 | 5.0 | 20 | 100 | 400 |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether | 138 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | 153 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | 154 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4',5-Hexabromodiphenyl ether | 156 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',3,4,4',5,6'-Heptabromodiphenyl ether | 183 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',3,4,4',6,6'-Heptabromodiphenyl ether | 184 | 2.0 | 10 | 40 | 200 | 800 |
| 2,3,3',4,4',5,6'-Heptabromodiphenyl ether | 191 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | 196 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | 197 | 2.0 | 10 | 40 | 200 | 800 |
| 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | 206 | 5.0 | 25 | 100 | 500 | 2000 |
| 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | 207 | 5.0 | 25 | 100 | 500 | 2000 |
| Decabromodiphenyl ether | 209 | 5.0 | 25 | 100 | 500 | 2000 |
| MASS-LABELLED PBDE CONGENERS | | | | | | |
| 4-Bromo[¹³ C ₁₂]diphenyl ether | 3L | 100 | 100 | 100 | 100 | 100 |
| 4,4'-Dibromo[¹³ C ₁₂]diphenyl ether | 15L | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 28L | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 47L | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 99L | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',6-Pentabromo[¹³ C ₁₂]diphenyl ether | 100L | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 126L | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 153L | 200 | 200 | 200 | 200 | 200 |
| 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 154L | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,4,4',5,6'-Heptabromo[¹³ C ₁₂]diphenyl ether | 183L | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,3',4,4',6,6'-Octabromo[¹³ C ₁₂]diphenyl ether | 197L | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,3',4,4',5,6,6'-Nonabromo[¹³ C ₁₂]diphenyl ether | 207L | 500 | 500 | 500 | 500 | 500 |
| Decabromo[¹³ C ₁₂]diphenyl ether | 209L | 500 | 500 | 500 | 500 | 500 |
| MASS-LABELLED PBDE INTERNAL STANDARD | | | | | | |
| 3,3',4,5'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 79L | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 138L | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,3',4,4',5,5'-Nonabromo[¹³ C ₁₂]diphenyl ether | 206L | 500 | 500 | 500 | 500 | 500 |

| Catalogue Number | Product (nonane/toluene solution) | Qty/Conc |
|-------------------|---|----------|
| MBDE-MXG | Mass-Labelled PBDE Solution/Mixture | 1.2 ml |
| MBDE-ISS-G | Mass-Labelled PBDE Internal Standard Solution | 1.2 ml |
| BDE-MXE | Native PBDE Solution/Mixture | 1.2 ml |

| NATIVE PBDE CONGENERS | IUPAC | MBDE-MXG (ng/ml) | MBDE-ISS-G (ng/ml) | BDE-MXE (ng/ml) |
|---|-------|---------------------|-----------------------|--------------------|
| 4-Bromodiphenyl ether | 3 | — | — | 1000 |
| 2,4-Dibromodiphenyl ether | 7 | — | — | 1000 |
| 4,4'-Dibromodiphenyl ether | 15 | — | — | 1000 |
| 2,2',4'-Tribromodiphenyl ether | 17 | — | — | 1000 |
| 2,4,4'-Tribromodiphenyl ether | 28 | — | — | 1000 |
| 2,2',4,4'-Tetrabromodiphenyl ether | 47 | — | — | 1000 |
| 2,2',4,5'-Tetrabromodiphenyl ether | 49 | — | — | 1000 |
| 2,3',4,4'-Tetrabromodiphenyl ether | 66 | — | — | 1000 |
| 2,3',4',6-Tetrabromodiphenyl ether | 71 | — | — | 1000 |
| 3,3',4,4'-Tetrabromodiphenyl ether | 77 | — | — | 1000 |
| 2,2',3,4,4'-Pentabromodiphenyl ether | 85 | — | — | 1000 |
| 2,2',4,4',5-Pentabromodiphenyl ether | 99 | — | — | 1000 |
| 2,2',4,4',6-Pentabromodiphenyl ether | 100 | — | — | 1000 |
| 2,3',4,4',6-Pentabromodiphenyl ether | 119 | — | — | 1000 |
| 3,3',4,4',5-Pentabromodiphenyl ether | 126 | — | — | 1000 |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether | 138 | — | — | 2000 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | 153 | — | — | 2000 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | 154 | — | — | 2000 |
| 2,3,3',4,4',5-Hexabromodiphenyl ether | 156 | — | — | 2000 |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | 183 | — | — | 2000 |
| 2,2',3,4,4',6,6'-Heptabromodiphenyl ether | 184 | — | — | 2000 |
| 2,3,3',4,4',5',6-Heptabromodiphenyl ether | 191 | — | — | 2000 |
| 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | 196 | — | — | 2000 |
| 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | 197 | — | — | 2000 |
| 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | 206 | — | — | 5000 |
| 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | 207 | — | — | 5000 |
| Decabromodiphenyl ether | 209 | — | — | 5000 |

MASS-LABELLED PBDE CONGENERS

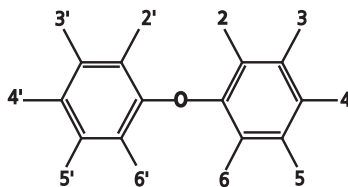
| | | | | |
|--|------|-----|---|---|
| 4-Bromo[¹³ C ₁₂]diphenyl ether | 3L | 100 | — | — |
| 4,4'-Dibromo[¹³ C ₁₂]diphenyl ether | 15L | 100 | — | — |
| 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 28L | 100 | — | — |
| 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 47L | 100 | — | — |
| 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 99L | 100 | — | — |
| 2,2',4,4',6-Pentabromo[¹³ C ₁₂]diphenyl ether | 100L | 100 | — | — |
| 3,3',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 126L | 100 | — | — |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 153L | 200 | — | — |
| 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 154L | 200 | — | — |
| 2,2',3,4,4',5',6-Heptabromo[¹³ C ₁₂]diphenyl ether | 183L | 200 | — | — |
| 2,2',3,3',4,4',6,6'-Octabromo[¹³ C ₁₂]diphenyl ether | 197L | 200 | — | — |
| 2,2',3,3',4,4',5,6,6'-Nonabromo[¹³ C ₁₂]diphenyl ether | 207L | 500 | — | — |
| Decabromo[¹³ C ₁₂]diphenyl ether | 209L | 500 | — | — |

MASS-LABELLED PBDE INTERNAL STANDARD

| | | | | |
|--|------|---|-----|---|
| 3,3',4,5'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 79L | — | 100 | — |
| 2,2',3,4,4',5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 138L | — | 200 | — |
| 2,2',3,3',4,4',5,5',6-Nonabromo[¹³ C ₁₂]diphenyl ether | 206L | — | 500 | — |

NATIVE BROMINATED DIPHENYL ETHERS (PBDEs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------|
| BDE-1 | 2-Bromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-2 | 3-Bromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-3 | 4-Bromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-7 | 2,4-Dibromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-10 | 2,6-Dibromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-15 | 4,4'-Dibromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-17 | 2,2',4-Tribromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-21 | 2,3,4-Tribromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-28 | 2,4,4'-Tribromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-30 | 2,4,6-Tribromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-37 | 3,4,4'-Tribromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-47 | 2,2',4,4'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-49 | 2,2',4,5'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-54 | 2,2',6,6'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-60 | 2,3,4,4'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-66 | 2,3',4,4'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-71 | 2,3',4',6-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-77 | 3,3',4,4'-Tetrabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-82 | 2,2',3,3',4-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-85 | 2,2',3,4,4'-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-99 | 2,2',4,4',5-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-100 | 2,2',4,4',6-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-104 | 2,2',4,6,6'-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-105 | 2,3,3',4,4'-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-119 | 2,3',4,4',6-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-126 | 3,3',4,4',5-Pentabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-128 | 2,2',3,3',4,4'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-138 | 2,2',3,4,4',5'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-139 | 2,2',3,4,4',6-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-140 | 2,2',3,4,4',6'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-149 | 2,2',3,4',5,6-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-153 | 2,2',4,4',5,5'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-154 | 2,2',4,4',5,6'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-155 | 2,2',4,4',6,6'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-156 | 2,3,3',4,4',5-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |

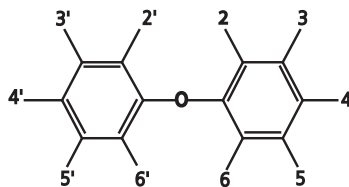


NATIVE BROMINATED DIPHENYL ETHERS (PBDEs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---------------------|---|-----------------|
| BDE-169 | 3,3',4,4',5,5'-Hexabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-170 | 2,2',3,3',4,4',5-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-171 | 2,2',3,3',4,4',6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-175 | 2,2',3,3',4,5',6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-176 | 2,2',3,3',4,6,6'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-177 | 2,2',3,3',4',5,6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-179 | 2,2',3,3',5,6,6'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-180 | 2,2',3,4,4',5,5'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-181 | 2,2',3,4,4',5,6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-182 | 2,2',3,4,4',5,6'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-183 | 2,2',3,4,4',5',6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-184 | 2,2',3,4,4',6,6'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-188 | 2,2',3,4',5,6,6'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-189 | 2,3,3',4,4',5,5'-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-191* | 2,3,3',4,4',5',6-Heptabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-194* | 2,2',3,3',4,4',5,5'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-195* | 2,2',3,3',4,4',5,6-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-196* | 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-197* | 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-198* | 2,2',3,3',4,5,5',6,-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-199* | 2,2',3,3',4,5,5',6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-200* | 2,2',3,3',4,5,6,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-201* | 2,2',3,3',4,5',6,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-202* | 2,2',3,3',5,5',6,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-203* | 2,2',3,4,4',5,5',6-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-204* | 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-205* | 2,3,3',4,4',5,5',6-Octabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-206* | 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-207* | 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-208* | 2,2',3,3',4,5,5',6,6'-Nonabromodiphenyl ether | 1.2 ml 50 µg/ml |
| BDE-209* | Decabromodiphenyl ether | 1.2 ml 50 µg/ml |
| 4PC-BDE-208* | 2,2',3,3',4,5,5',6,6'-Nonabromo-4'-chlorodiphenyl ether | 1.2 ml 50 µg/ml |

4PC-BDE-208 may be useful as an internal or surrogate standard for HRGC/ECD, HRGC/FID, and/or HRGC/MS analyses.

* Toluene solution



PBDE WINDOW DEFINING SOLUTION/MIXTURE

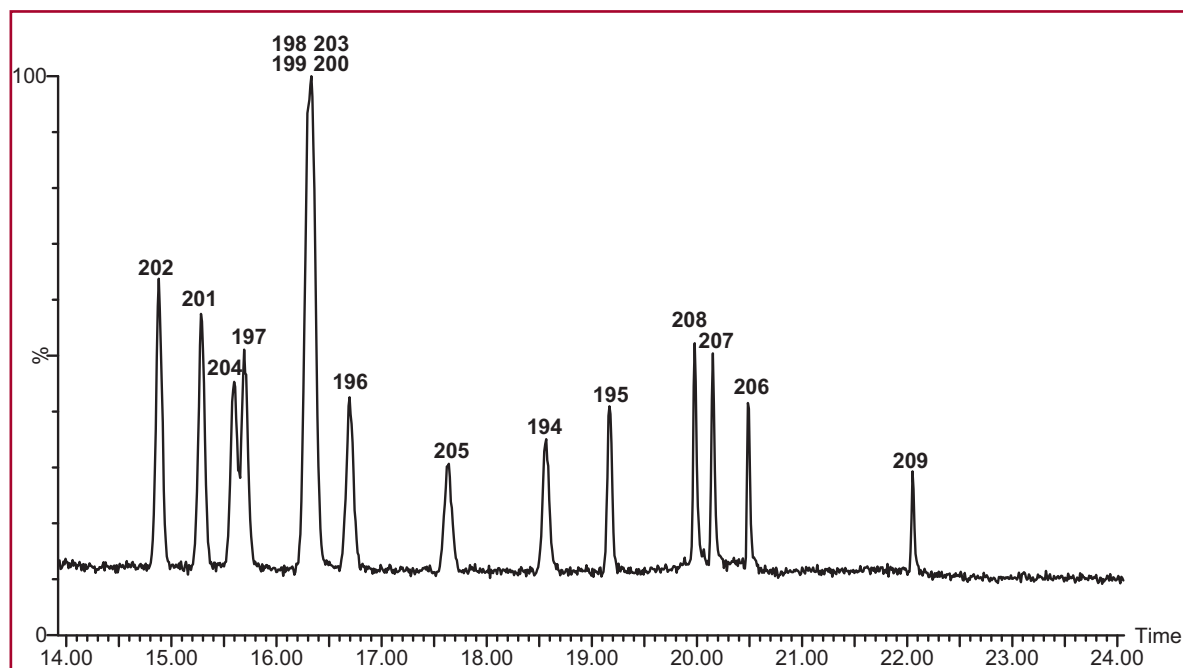
| Catalogue Number | Product (nonane solution) | Qty/Conc | |
|---------------------------|---|---------------------------------|---------------------------------------|
| BDE-WD | PBDE Window Defining Solution/Mixture for use with a J&W DB-5HT column | 1.2 ml | |
| | | FIRST ELUTER (IUPAC) | LAST ELUTER (IUPAC) |
| | | | CONCENTRATION (µg/ml each) |
| Bromodiphenyl ethers | | 1 | 3 |
| Dibromodiphenyl ethers | | 10 | 15 |
| Tribromodiphenyl ethers | | 30 | 37 |
| Tetrabromodiphenyl ethers | | 54 | 60 |
| Pentabromodiphenyl ethers | | 104 | 82 |
| Hexabromodiphenyl ethers | | 155 | 128 |
| Heptabromodiphenyl ethers | | 188 | 170 |
| Octabromodiphenyl ethers | | 202 | 195 |
| Nonabromodiphenyl ethers | | 208 | 206 |
| Decabromodiphenyl ether | | | 209 |
| | | | 5.0 |
| | | | 5.0 |

NATIVE BROMINATED DIPHENYL ETHERS: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc | |
|---|------------------------------|--------------|----------|
| BDE-MXA | Native PBDE Solution/Mixture | 1.2 ml | |
| | | IUPAC | |
| 2,2',4,4'-Tetrabromodiphenyl ether | | 47 | 5 µg/ml |
| 2,2',4,4',5-Pentabromodiphenyl ether | | 99 | 5 µg/ml |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | | 153 | 5 µg/ml |
| BDE-MXB | Native PBDE Solution/Mixture | 1.2 ml | |
| | | IUPAC | |
| 2,4,4'-Tribromodiphenyl ether | | 28 | 5 µg/ml |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether | | 154 | 5 µg/ml |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | | 183 | 5 µg/ml |
| BDE-MXD | Native PBDE Solution/Mixture | 1.2 ml | |
| | | IUPAC | |
| 2,2',4-Tribromodiphenyl ether | | 17 | 5 µg/ml |
| 2,2',4,4'-Tetrabromodiphenyl ether | | 47 | 5 µg/ml |
| 2,3',4,4'-Tetrabromodiphenyl ether | | 66 | 5 µg/ml |
| 2,2',4,4',6-Pentabromodiphenyl ether | | 100 | 5 µg/ml |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether | | 153 | 5 µg/ml |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether | | 183 | 5 µg/ml |
| Decabromodiphenyl ether | | 209 | 10 µg/ml |

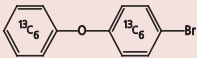
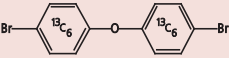
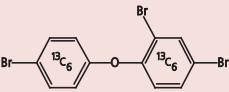
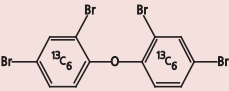
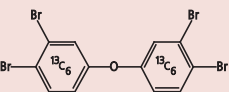
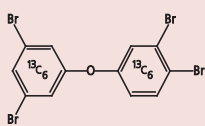
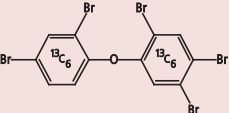
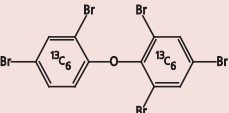
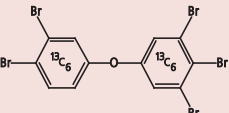
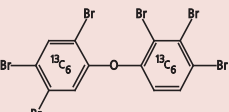
NATIVE BROMINATED DIPHENYL ETHERS: SOLUTION/MIXTURE

| Catalogue Number | Product (nonane/toluene solution) | Qty/Conc |
|---|---|-----------|
| BDE-OND | Solution/Mixture of Octa-, Nona-, and Deca-BDEs | 1.2 ml |
| | IUPAC | |
| 2,2',3,3',4,4',5,5'-Octabromodiphenyl ether | 194 | 1.0 µg/ml |
| 2,2',3,3',4,4',5,6-Octabromodiphenyl ether | 195 | 1.0 µg/ml |
| 2,2',3,3',4,4',5,6'-Octabromodiphenyl ether | 196 | 1.0 µg/ml |
| 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | 197 | 1.0 µg/ml |
| 2,2',3,3',4,5,5',6-Octabromodiphenyl ether | 198 | 1.0 µg/ml |
| 2,2',3,3',4,5,5',6'-Octabromodiphenyl ether | 199 | 1.0 µg/ml |
| 2,2',3,3',4,5,6,6'-Octabromodiphenyl ether | 200 | 1.0 µg/ml |
| 2,2',3,3',4,5',6,6'-Octabromodiphenyl ether | 201 | 1.0 µg/ml |
| 2,2',3,3',5,5',6,6'-Octabromodiphenyl ether | 202 | 1.0 µg/ml |
| 2,2',3,4,4',5,5',6-Octabromodiphenyl ether | 203 | 1.0 µg/ml |
| 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether | 204 | 1.0 µg/ml |
| 2,3,3',4,4',5,5',6-Octabromodiphenyl ether | 205 | 1.0 µg/ml |
| 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | 206 | 2.5 µg/ml |
| 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | 207 | 2.5 µg/ml |
| 2,2',3,3',4,5,5',6,6'-Nonabromodiphenyl ether | 208 | 2.5 µg/ml |
| Decabromodiphenyl ether | 209 | 2.5 µg/ml |



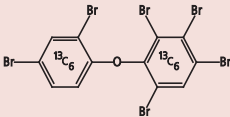
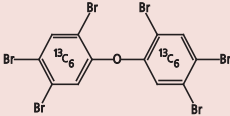
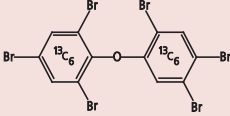
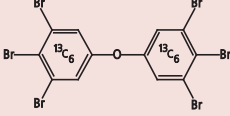
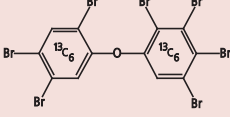
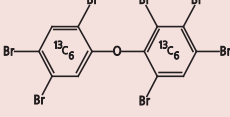
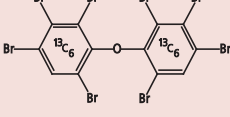
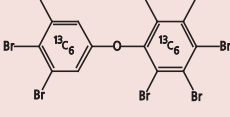
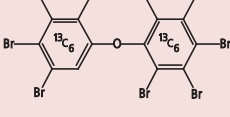
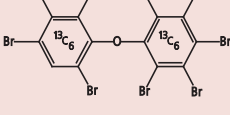
HRGC/HRMS Data for BDE-OND on a 15m DB-5HT column.

MASS-LABELLED BROMINATED DIPHENYL ETHERS

| Catalogue Number | Product |
|------------------|---|
| MBDE-3 |  <p>4-Bromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-15 |  <p>4,4'-Dibromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-28 |  <p>2,4,4'-Tribromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-47 |  <p>2,2',4,4'-Tetrabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-77 |  <p>3,3',4,4'-Tetrabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-79 |  <p>3,3',4,5'-Tetrabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-99 |  <p>2,2',4,4',5-Pentabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-100 |  <p>2,2',4,4',6-Pentabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-126 |  <p>3,3',4,4',5-Pentabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-138 |  <p>2,2',3,4,4',5'-Hexabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

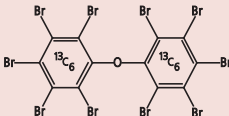
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED BROMINATED DIPHENYL ETHERS

| Catalogue Number | Product |
|------------------|---|
| MBDE-139 |  <p>2,2',3,4,4',6-Hexabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-153 |  <p>2,2',4,4',5,5'-Hexabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-154 |  <p>2,2',4,4',5,6'-Hexabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-169 |  <p>3,3',4,4',5,5'-Hexabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-180 |  <p>2,2',3,4,4',5,5'-Heptabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-183 |  <p>2,2',3,4,4',5,6-Heptabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MBDE-197 |  <p>2,2',3,3',4,4',6,6'-Octabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MBDE-205 |  <p>2,3,3',4,4',5,5',6-Octabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MBDE-206 |  <p>2,2',3,3',4,4',5,5',6-Nonabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| MBDE-207 |  <p>2,2',3,3',4,4',5,6,6'-Nonabromo[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED BROMINATED DIPHENYL ETHERS

| Catalogue Number | Product |
|------------------|---|
| MBDE-209 |  <p>Decabromo[¹³C₁₂]diphenyl ether 1.2 ml; 25 µg/ml (±1.2 µg/ml); in toluene</p> |

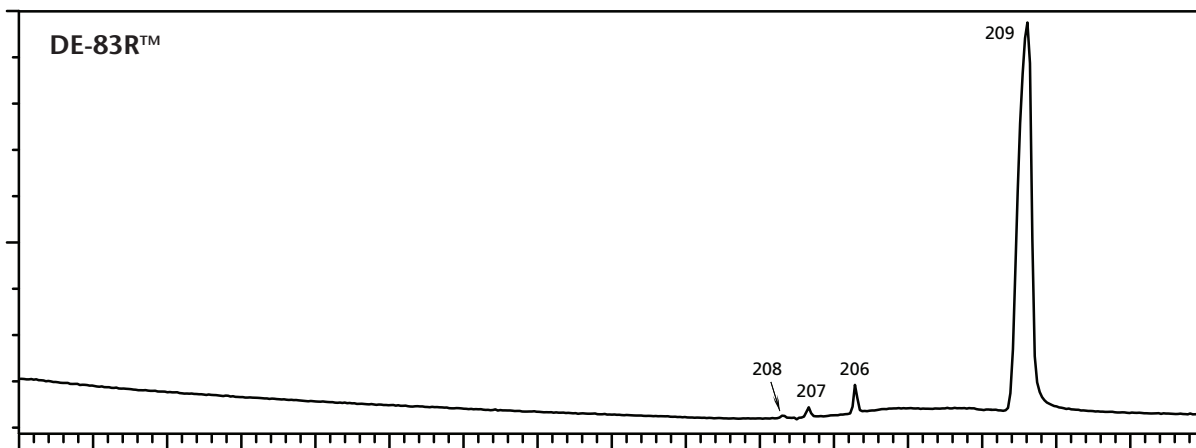
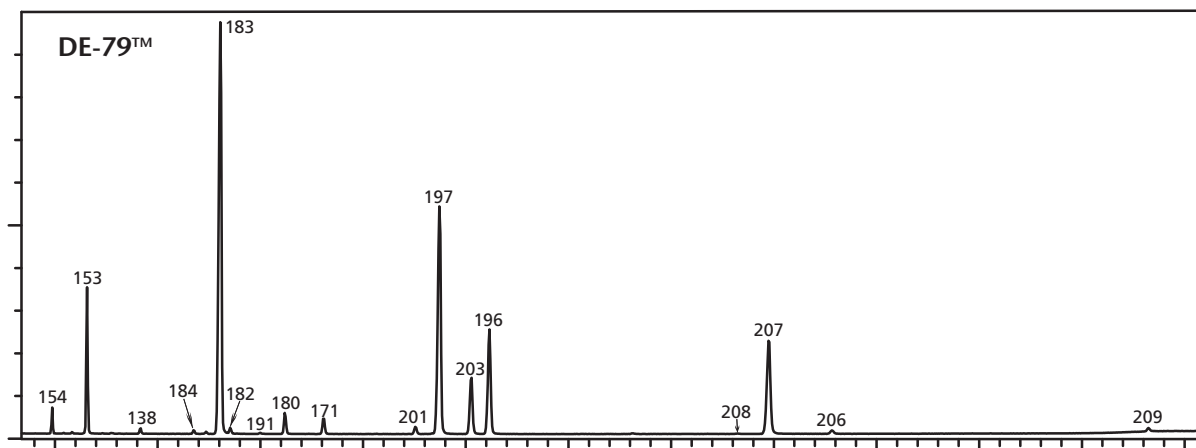
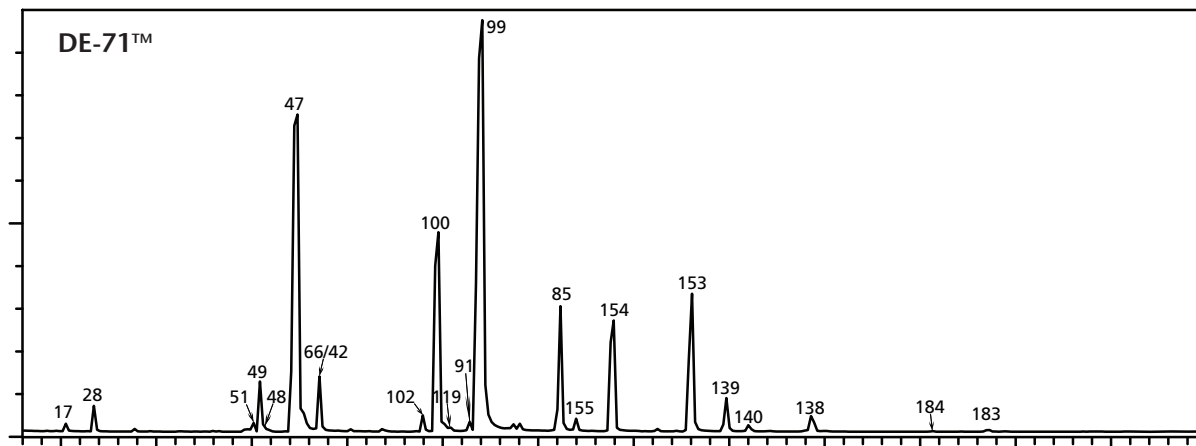
* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED BROMINATED DIPHENYL ETHERS: SOLUTION/MIXTURES

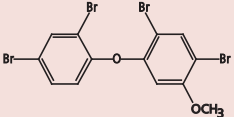
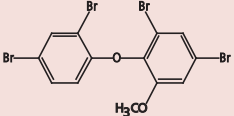
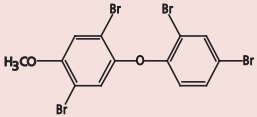
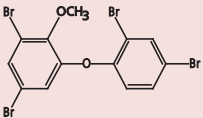
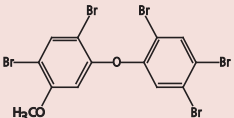
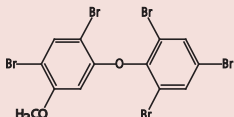
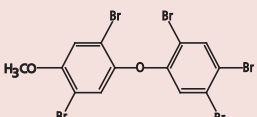
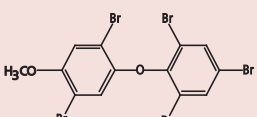
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|----------|
| MBDE-MXA | Mass-Labelled PBDE Solution/Mixture | 1.2 ml |
| | IUPAC | |
| | 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| MBDE-MXB | Mass-Labelled PBDE Solution/Mixture | 1.2 ml |
| | IUPAC | |
| | 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',3,4,4',5',6-Heptabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| MBDE-MXC | Mass-Labelled PBDE Solution/Mixture | 1.2 ml |
| | IUPAC | |
| | 4-Bromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 4,4'-Dibromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,4,4'-Tribromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4'-Tetrabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5-Pentabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',4,4',5,6'-Hexabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |
| | 2,2',3,4,4',5',6-Heptabromo[¹³ C ₁₂]diphenyl ether | 5 µg/ml |

BROMINATED DIPHENYL ETHER TECHNICAL MIXTURES

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|------------------|
| TBDE-71 | Great Lakes Chemical DE-71™ Pentabromodiphenyl Oxide | 1.2 ml 100 µg/ml |
| TBDE-79 | Great Lakes Chemical DE-79™ Octabromodiphenyl Oxide | 1.2 ml 100 µg/ml |
| TBDE-83R | Great Lakes Chemical DE-83R™ Decabromodiphenyl Oxide | 1.2 ml 100 µg/ml |



NATIVE METHOXY-BROMODIPHENYL ETHERS (MeO-BDEs)

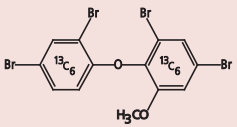
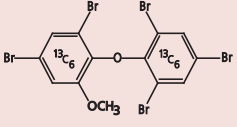
| Catalogue Number | Product |
|------------------|---|
| 5MBDE47 |  <p>2,2',4,4'-Tetrabromo-5-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 6MBDE47 |  <p>2,2',4,4'-Tetrabromo-6-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 4PMBDE49 |  <p>2,2',4',5-Tetrabromo-4-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 2PMBDE68 |  <p>2',3,4',5-Tetrabromo-2-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 5PMBDE99 |  <p>2,2',4,4',5-Pentabromo-5'-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 5PMBDE100 |  <p>2,2',4,4',6'-Pentabromo-5-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 4PMBDE101 |  <p>2,2',4,5,5'-Pentabromo-4'-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| 4PMBDE103 |  <p>2,2',4',5,6'-Pentabromo-4-methoxydiphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

NOTE: In order to emphasize their relationship to the corresponding PBDE, these compounds have been numbered based on the diphenyl ether as the parent molecule with bromines retaining their BDE numbers. The methoxy groups are treated as additional substituents and listed alphabetically.

METHOXY-BROMODIPHENYL ETHERS: SOLUTION/MIXTURE

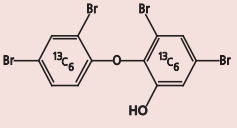
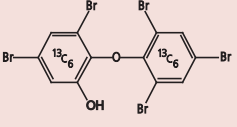
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| MeOBDES | Methoxy-Bromodiphenyl Ethers Solution/Mixture | 1.2 ml |
| | 2,2',4,4'-Tetrabromo-5-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4,4'-Tetrabromo-6-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4',5-Tetrabromo-4-methoxydiphenyl ether | 5 µg/ml |
| | 2',3,4',5-Tetrabromo-2-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4,4',5-Pentabromo-5'-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4,4',6'-Pentabromo-5-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4,5,5'-Pentabromo-4'-methoxydiphenyl ether | 5 µg/ml |
| | 2,2',4',5,6'-Pentabromo-4-methoxydiphenyl ether | 5 µg/ml |

MASS-LABELLED METHOXY-BROMODIPHENYL ETHERS

| Catalogue Number | Product |
|-------------------|---|
| M6MBDE47 |  <p>2,2',4,4'-Tetrabromo-6-methoxy[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| M6PMBDE100 |  <p>2,2',4,4',6-Pentabromo-6'-methoxy[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED HYDROXY-BROMODIPHENYL ETHERS

| Catalogue Number | Product |
|-------------------|---|
| M6HBDE47 |  <p>2,2',4,4'-Tetrabromo-6-hydroxy[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M6PHBDE100 |  <p>2,2',4,4',6-Pentabromo-6'-hydroxy[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

NOTE: In order to emphasize their relationship to the corresponding PBDE, these compounds have been numbered based on the diphenyl ether as the parent molecule with bromines retaining their BDE numbers. The methoxy and hydroxy groups are treated as additional substituents and listed alphabetically.

NATIVE BROMINATED BIPHENYLS (PBBs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-----------------|
| BB-1 | 2-Bromobiphenyl | 1.2 ml 50 µg/ml |
| BB-2 | 3-Bromobiphenyl | 1.2 ml 50 µg/ml |
| BB-3 | 4-Bromobiphenyl | 1.2 ml 50 µg/ml |
| BB-4 | 2,2'-Dibromobiphenyl | 1.2 ml 50 µg/ml |
| BB-7 | 2,4-Dibromobiphenyl | 1.2 ml 50 µg/ml |
| BB-9 | 2,5-Dibromobiphenyl | 1.2 ml 50 µg/ml |
| BB-10 | 2,6-Dibromobiphenyl | 1.2 ml 50 µg/ml |
| BB-15 | 4,4'-Dibromobiphenyl | 1.2 ml 50 µg/ml |
| BB-18 | 2,2',5-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-22 | 2,3,4'-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-26 | 2,3',5-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-29 | 2,4,5-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-30 | 2,4,6-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-31 | 2,4',5-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-37 | 3,4,4'-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-38 | 3,4,5-Tribromobiphenyl | 1.2 ml 50 µg/ml |
| BB-49 | 2,2',4,5'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-52 | 2,2',5,5'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-53 | 2,2',5,6'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-56 | 2,3,3',4'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-75 | 2,4,4',6-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-77 | 3,3',4,4'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-80 | 3,3',5,5'-Tetrabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-101 | 2,2',4,5,5'-Pentabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-103 | 2,2',4,5',6-Pentabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-153 | 2,2',4,4',5,5'-Hexabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-154 | 2,2',4,4',5,6'-Hexabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-155 | 2,2',4,4',6,6'-Hexabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-156 | 2,3,3',4,4',5-Hexabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-169 | 3,3',4,4',5,5'-Hexabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-180 | 2,2',3,4,4',5,5'-Heptabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-194 | 2,2',3,3',4,4',5,5'-Octabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-205 | 2,3,3',4,4',5,5',6-Octabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-206* | 2,2',3,3',4,4',5,5',6-Nonabromobiphenyl | 1.2 ml 50 µg/ml |
| BB-209* | Decabromobiphenyl | 1.2 ml 50 µg/ml |

* 50% Nonane/50% Toluene Solution

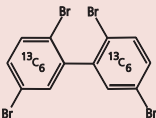
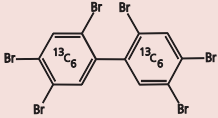
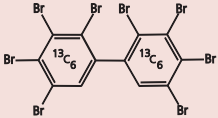
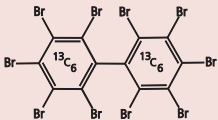
NATIVE BROMINATED BIPHENYLS: SOLUTION/MIXTURE

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|-----------------------------|----------|
| PBB-MXA | Native PBB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 4-Bromobiphenyl | 3 | 1 µg/ml |
| 4,4'-Dibromobiphenyl | 15 | 1 µg/ml |
| 2,2',5-Tribromobiphenyl | 18 | 1 µg/ml |
| 2,2',5,5'-Tetrabromobiphenyl | 52 | 1 µg/ml |
| 2,2',4,5,5'-Pentabromobiphenyl | 101 | 2 µg/ml |
| 2,2',4,4',5,5'-Hexabromobiphenyl | 153 | 2 µg/ml |
| 2,2',3,4,4',5,5'-Heptabromobiphenyl | 180 | 2 µg/ml |
| 2,2',3,3',4,4',5,5'-Octabromobiphenyl | 194 | 2 µg/ml |
| 2,2',3,3',4,4',5,5',6-Nonabromobiphenyl | 206 | 5 µg/ml |
| Decabromobiphenyl | 209 | 5 µg/ml |

POLYBROMINATED BIPHENYL TECHNICAL MIXTURES

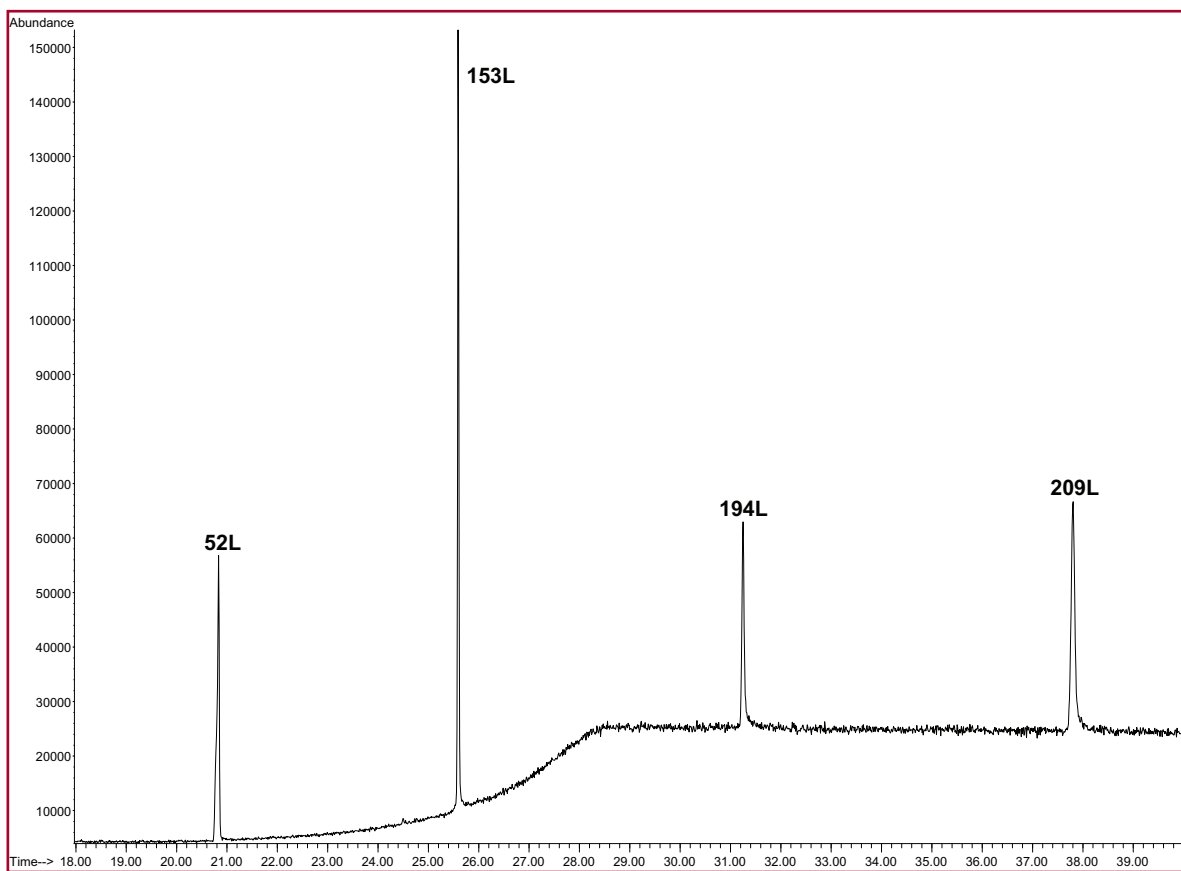
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|------------------|
| TBB-BP6 | Great Lakes Chemical Firemaster BP-6™ Hexabromobiphenyl | 1.2 ml 100 µg/ml |
| TBB-809D | Chemische Fabrik Kalk Bromkal80-9D™ Nonabromobiphenyl | 1.2 ml 100 µg/ml |

MASS-LABELLED BROMINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|---|
| MBB-52 |  <p>2,2',5,5'-Tetrabromo[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic purity; 99% or greater</p> |
| MBB-153 |  <p>2,2',4,4',5,5'-Hexabromo[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic purity; 99% or greater</p> |
| MBB-194 |  <p>2,2',3,3',4,4',5,5'-Octabromo[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic purity; 99% or greater</p> |
| MBB-209 |  <p>Decabromo[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic purity; 99% or greater</p> |

MASS-LABELLED BROMINATED BIPHENYLS: SOLUTION/MIXTURE

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--|------------------------------------|----------|
| MBB-MXA | Mass-Labelled PBB Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 2,2',5,5'-Tetrabromo[¹³ C ₁₂]biphenyl | 52L | 1 µg/ml |
| 2,2',4,4',5,5'-Hexabromo[¹³ C ₁₂]biphenyl | 153L | 2 µg/ml |
| 2,2',3,3',4,4',5,5'-Octabromo[¹³ C ₁₂]biphenyl | 194L | 2 µg/ml |
| Decabromo[¹³ C ₁₂]biphenyl | 209L | 5 µg/ml |



HRGC/LRMS Data for MBB-MXA on a 30m DB-5 column.



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HALOGENATED FLAME RETARDANTS & RELATED COMPOUNDS

This section is devoted to halogenated flame retardants (HFRs) and related compounds. Since 2012, we have added a number of flame retardants, most notably:

Native and mass-labelled Hexabromocyclododecane (HBCD) mixtures

Native Experimental Flame Retardants (EFRs)

Native and mass-labelled Organophosphorus Compounds

2-Ethylhexyl-d₁₇-2,3,4,5-tetrabromo[¹³C₆]benzoate (MEHTBB)

Bis(2-Ethylhexyl-d₁₇)-tetrabromo[¹³C₆]phthalate (MBEHTBP)

This section also contains compounds related to HFRs that are created through metabolism, combustion or other processes, namely:

Native and mass-labelled PBDDs

Native PBDFs

Mixed Br/Cl Dioxins and Furans

Native and mass-labelled Bromophenols

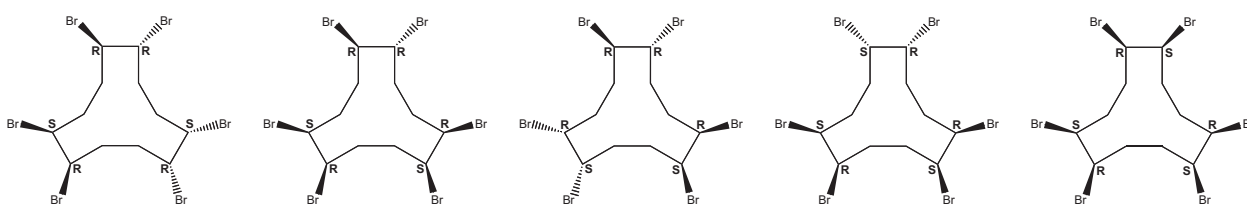


NATIVE HEXABROMOCYCLODODECANE ISOMERS (HBCDs)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|----------------------|
| aHBCD | α -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| bHBCD | β -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| gHBCD | γ -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| dHBCD | δ -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| eHBCD | ε -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| zHBCD | ζ -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| etaHBCD | η -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| tHBCD | θ -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| iHBCD | ι -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |
| kHBCD | κ -1,2,5,6,9,10-Hexabromocyclododecane | 1.2 ml 50 μ g/ml |

NATIVE HBCD ISOMERS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|---------------|
| HBCD-MXA | Native Hexabromocyclododecane Isomer Solution/Mixture | 1.2 ml |
| | α -1,2,5,6,9,10-Hexabromocyclododecane | 10 μ g/ml |
| | β -1,2,5,6,9,10-Hexabromocyclododecane | 10 μ g/ml |
| | γ -1,2,5,6,9,10-Hexabromocyclododecane | 10 μ g/ml |



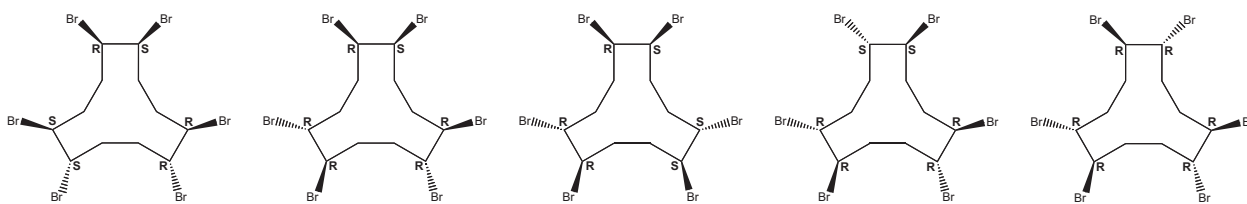
alpha

beta

gamma

delta

epsilon



zeta

eta

theta

iota

kappa

¹³C-LABELLED HEXABROMOCYCLODODECANE ISOMERS

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|-----------------|
| MaHBCD | α -1,2,5,6,9,10-Hexabromo[¹³ C ₁₂]cyclododecane | 1.2 ml 50 µg/ml |
| MbHBCD | β -1,2,5,6,9,10-Hexabromo[¹³ C ₁₂]cyclododecane | 1.2 ml 50 µg/ml |
| MgHBCD | γ -1,2,5,6,9,10-Hexabromo[¹³ C ₁₂]cyclododecane | 1.2 ml 50 µg/ml |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

DEUTERATED HEXABROMOCYCLODODECANE ISOMERS

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|-----------------|
| DaHBCD | α -1,2,5,6,9,10-Hexabromocyclododecane-d ₁₈ | 1.2 ml 50 µg/ml |
| DbHBCD | β -1,2,5,6,9,10-Hexabromocyclododecane-d ₁₈ | 1.2 ml 50 µg/ml |
| DgHBCD | γ -1,2,5,6,9,10-Hexabromocyclododecane-d ₁₈ | 1.2 ml 50 µg/ml |

* Unless stated otherwise, isotopic purities of these compounds are 98%.

¹³C-LABELLED HBCD ISOMERS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|----------|
| MHBCD-MXA | Mass-Labelled Hexabromocyclododecane Isomer Solution/Mixture | 1.2 ml |
| | α -1,2,5,6,9,10-Hexabromocyclo[¹³ C ₁₂]dodecane | 10 µg/ml |
| | β -1,2,5,6,9,10-Hexabromocyclo[¹³ C ₁₂]dodecane | 10 µg/ml |
| | γ -1,2,5,6,9,10-Hexabromocyclo[¹³ C ₁₂]dodecane | 10 µg/ml |

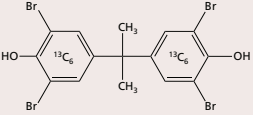
PENTABROMOCYCLODODECENE (PBCD)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|-----------------|
| PBCD | <i>rac</i> -(1,5 <i>R</i> ,6,5,9 <i>S</i> ,10 <i>R</i>)-pentabromocyclododecene | 1.2 ml 50 µg/ml |

TETRABROMOBISPHENOL-A (TBBPA)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---------------------------------|-----------------|
| TBBPA | 3,3',5,5'-Tetrabromobisphenol-A | 1.2 ml 50 µg/ml |

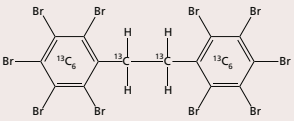
MASS-LABELLED TETRABROMOBISPHENOL-A

| Catalogue Number | Product |
|------------------|---|
| MTBBPA |  <p>3,3',5,5'-Tetrabromobisphenol-A [rings; ¹³C₁₂] 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; Isotopic purity; 99% or greater</p> |

DECABROMODIPHENYLETHANE (DBDPE)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---------------------------------|-----------------|
| DBDPE | 1,2-Bis(pentabromophenyl)ethane | 1.2 ml 25 µg/ml |

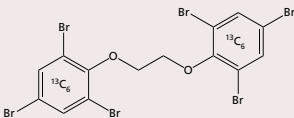
MASS-LABELLED DECABROMODIPHENYLETHANE

| Catalogue Number | Product |
|------------------|---|
| MDBDPE |  <p>1,2-Bis(pentabromophenyl)ethane [¹³C₁₄] 1.2 ml; 25 µg/ml (±1.2 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |

1,2-BIS(2,4,6-TRIBROMOPHENOXY)ETHANE (BTBPE)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--------------------------------------|-----------------|
| BTBPE | 1,2-Bis(2,4,6-tribromophenoxy)ethane | 1.2 ml 50 µg/ml |

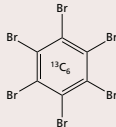
MASS-LABELLED 1,2-BIS(2,4,6-TRIBROMOPHENOXY)ETHANE

| Catalogue Number | Product |
|------------------|---|
| MBTBPE |  <p>1,2-Bis(2,4,6-tribromo[¹³C₆]phenoxy)ethane 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic purity; 99% or greater</p> |

HEXABROMOBENZENE (HBBZ)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------|-----------------|
| HBBZ | Hexabromobenzene | 1.2 ml 50 µg/ml |

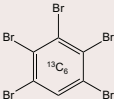
MASS-LABELLED HEXABROMOBENZENE

| Catalogue Number | Product |
|------------------|--|
| MHBBZ |  <p>Hexabromo[¹³C₆]benzene 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |

PENTABROMOBENZENE (PBBZ)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------|-----------------|
| PBBZ | Pentabromobenzene | 1.2 ml 50 µg/ml |

MASS-LABELLED PENTABROMOBENZENE

| Catalogue Number | Product |
|------------------|--|
| MPBBZ |  Pentabromo[¹³ C ₆]benzene 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater |

PENTABROMOETHYLBENZENE (PBEB)

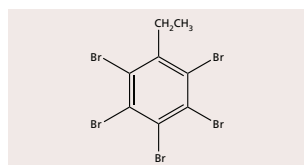
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------|-----------------|
| PBEB | Pentabromoethylbenzene | 1.2 ml 50 µg/ml |

PENTABROMOTOLUENE (PBT)

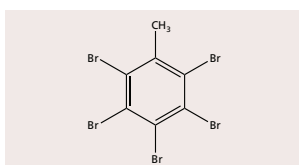
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------|-----------------|
| PBT | Pentabromotoluene | 1.2 ml 50 µg/ml |

TETRABROMO-P-XYLENE (pTBX)

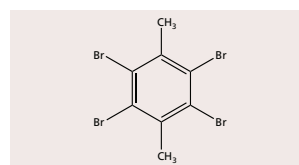
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|-----------------------------|-----------------|
| pTBX | 2,3,5,6-Tetrabromo-p-xylene | 1.2 ml 50 µg/ml |



Pentabromoethylbenzene



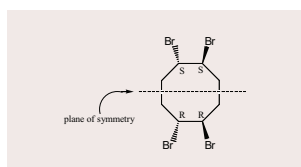
Pentabromotoluene



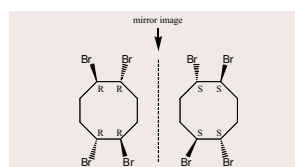
2,3,5,6-Tetrabromo-p-xylene

1,2,5,6-TETRABROMOCYCLOOCTANE (TBCO)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|-----------------|
| aTBCO | (1R,2R,5S,6S)-1,2,5,6-tetrabromocyclooctane | 1.2 ml 50 µg/ml |
| bTBCO | rac-(1R,2R,5R,6R)-1,2,5,6-tetrabromocyclooctane | 1.2 ml 50 µg/ml |



aTBCO



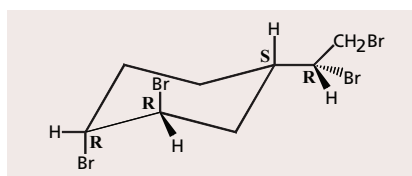
bTBCO

TETRABROMOETHYLCYCLOHEXANE (TBECH)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|-----------------|
| bTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>S</i>)-4-((1 <i>S</i>)-1,2-dibromoethyl)cyclohexane | 1.2 ml 50 µg/ml |
| gTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>R</i>)-4-((1 <i>R</i>)-1,2-dibromoethyl)cyclohexane | 1.2 ml 50 µg/ml |

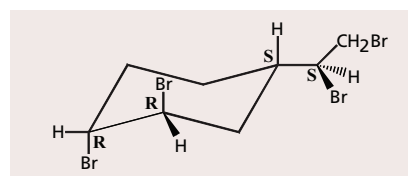
TETRABROMOETHYLCYCLOHEXANE ISOMERIC SOLUTION/MIXTURES

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|----------|
| abTBECH | TBECH Isomeric Solution/Mixture | 1.2 ml |
| aTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>S</i>)-4-((1 <i>R</i>)-1,2-dibromoethyl)cyclohexane | 50 µg/ml |
| bTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>S</i>)-4-((1 <i>S</i>)-1,2-dibromoethyl)cyclohexane | 50 µg/ml |
| gdTBECH | TBECH Isomeric Solution/Mixture | 1.2 ml |
| gTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>R</i>)-4-((1 <i>R</i>)-1,2-dibromoethyl)cyclohexane | 50 µg/ml |
| dTBECH | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i>)-1,2-dibromo-(4 <i>R</i>)-4-((1 <i>S</i>)-1,2-dibromoethyl)cyclohexane | 50 µg/ml |



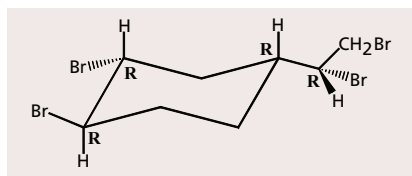
aTBECH

rac-(1*R*,2*R*)-1,2-dibromo-(4*S*)-4-((1*R*)-1,2-dibromoethyl)cyclohexane



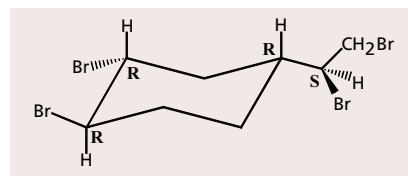
bTBECH

rac-(1*R*,2*R*)-1,2-dibromo-(4*S*)-4-((1*S*)-1,2-dibromoethyl)cyclohexane



gTBECH

rac-(1*R*,2*R*)-1,2-dibromo-(4*R*)-4-((1*R*)-1,2-dibromoethyl)cyclohexane

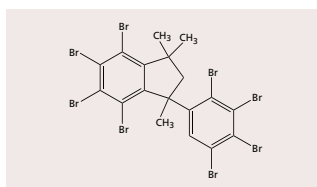


dTBECH

rac-(1*R*,2*R*)-1,2-dibromo-(4*R*)-4-((1*S*)-1,2-dibromoethyl)cyclohexane

OCTABROMOTRIMETHYLPHENYLINDANE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|-----------------|
| OBIND | 4,5,6,7-Tetrabromo-1,1,3-trimethyl-3-(2,3,4,5-tetrabromophenyl)-indane | 1.2 ml 50 µg/ml |

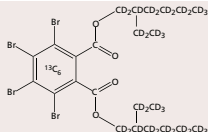


OBIND

BIS(2-ETHYLHEXYL)TETRABROMOPHTHALATE (BEHTBP)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|-----------------|
| BEHTBP | Bis(2-ethyl-1-hexyl)tetrabromophthalate | 1.2 ml 50 µg/ml |

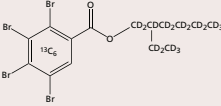
BIS(2-ETHYLHEXYL-d₁₇)TETRABROMO[¹³C₆]PHTHALATE (MBEHTBP)

| Catalogue Number | Product |
|------------------|---|
| MBEHTBP |  <p>Bis(2-ethylhexyl-d₁₇)tetrabromo[¹³C₆]phthalate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater (¹³C₆); 98% or greater (²H₃₄)</p> |

2-ETHYLHEXYL-2,3,4,5-TETRABROMOBENZOATE (EHTBB)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---|-----------------|
| EHTBB | 2-Ethylhexyl-2,3,4,5-tetrabromobenzoate | 1.2 ml 50 µg/ml |

2-ETHYLHEXYL-d₁₇-2,3,4,5-TETRABROMO[¹³C₆]BENZOATE (MEHTBB)

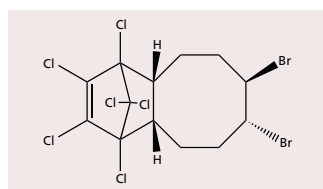
| Catalogue Number | Product |
|------------------|--|
| MEHTBB |  <p>2-Ethylhexyl-d₁₇-2,3,4,5-tetrabromo[¹³C₆]benzoate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater (¹³C₆); 98% or greater (²H₁₇)</p> |

HEXACHLOROCYCLOPENTENYL-DIBROMOCYCLOOCTANE (HCDBCO)

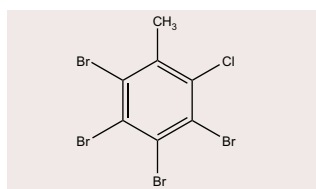
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|-----------------|
| HCDBCO | <i>rac</i> -(1 <i>R</i> ,2 <i>R</i> ,5 <i>R</i> ,6 <i>R</i> ,9 <i>S</i> ,10 <i>S</i>)-5,6-dibromo-1,10,11,12,13,13-hexachlorotricyclo[8.2.1.0 ^{2,9}]tridec-11-ene | 1.2 ml 50 µg/ml |

TETRABROMO-O-CHLOROTOLUENE (TBCT)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|-------------------------------------|-----------------|
| TBCT | Tetrabromo- <i>o</i> -chlorotoluene | 1.2 ml 50 µg/ml |



HCDBCO



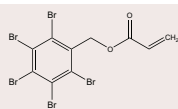
TBCT

PENTABROMOBENZYL ACRYLATE

Catalogue Number

Product

PBBA



Pentabromobenzyl acrylate

1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene

2,4,6-TRIBROMOPHENYL ETHERS

Catalogue Number

Product (toluene solution)

Qty/Conc

ATE

Allyl 2,4,6-tribromophenyl ether

1.2 ml 50 µg/ml

DPTE

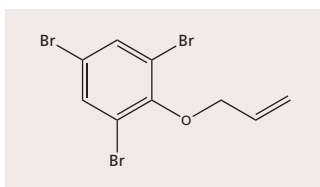
2,3-Dibromopropyl 2,4,6-tribromophenyl ether

1.2 ml 50 µg/ml

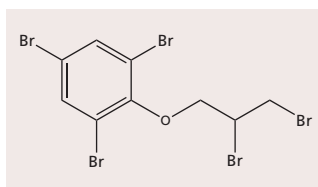
BATE

2-Bromoallyl 2,4,6-tribromophenyl ether

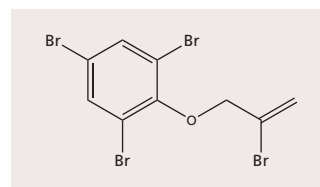
1.2 ml 50 µg/ml



ATE



DPTE



BATE

DECHLORANE PLUS®

Catalogue Number

Product (toluene solution)

Qty/Conc

s-DP

syn-Dechlorane Plus®

1.2 ml 50 µg/ml

a-DP

anti-Dechlorane Plus®

1.2 ml 50 µg/ml

DECHLORINATED DECHLORANE PLUS®

Catalogue Number

Product (toluene solution)

Qty/Conc

aCl10DP

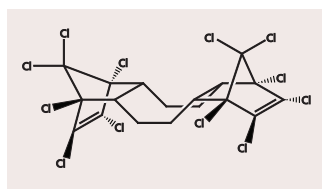
Cl10 Dechlorane Plus®

1.2 ml 50 µg/ml

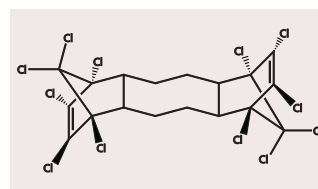
aCl11DP

Cl11 Dechlorane Plus®

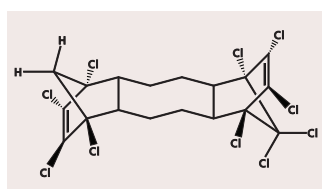
1.2 ml 50 µg/ml



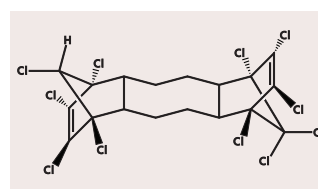
syn-Dechlorane Plus®



anti-Dechlorane Plus®



anti-Cl10 Dechlorane Plus®



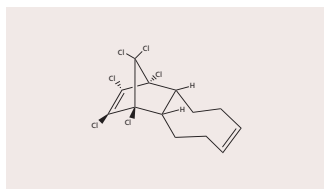
anti-Cl11 Dechlorane Plus®

DECHLORANE PLUS® MONO ADDUCTS

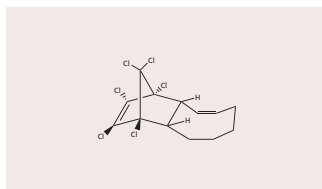
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------------|-----------------|
| DPMA | Dechlorane Plus® Mono adduct | 1.2 ml 50 µg/ml |
| 1,3-DPMA | 1,3-Dechlorane Plus® Mono adduct | 1.2 ml 50 µg/ml |

EXPERIMENTAL FLAME RETARDANTS (EFRs)

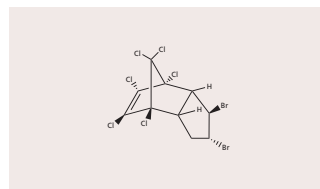
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|------------------------------|-----------------|
| DBCD | Dibromochlordene | 1.2 ml 50 µg/ml |
| Dec-601 | Dechlorane 601 | 1.2 ml 50 µg/ml |
| Dec-602 | Dechlorane 602 | 1.2 ml 50 µg/ml |
| Dec-603 | Dechlorane 603 | 1.2 ml 50 µg/ml |
| Dec-604 | Dechlorane 604 | 1.2 ml 50 µg/ml |
| Dec-604CB | Dechlorane 604 Component B | 1.2 ml 50 µg/ml |
| CPlus | Chlordene Plus | 1.2 ml 50 µg/ml |
| DBALD | Dibromoaldrin | 1.2 ml 50 µg/ml |
| HCPN | Hexachloro(phenyl)norbornene | 1.2 ml 50 µg/ml |



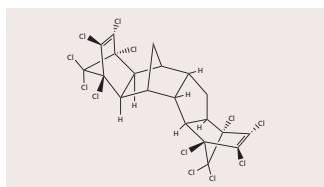
DPMA



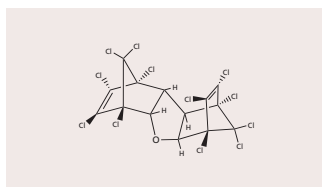
1,3-DPMA



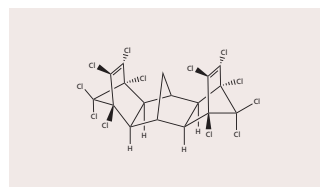
DBCD



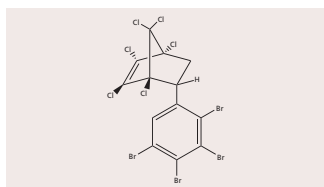
Dec-601



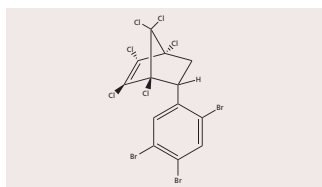
Dec-602



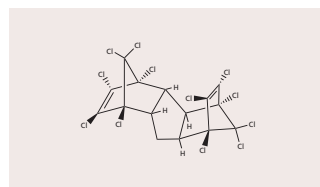
Dec-603



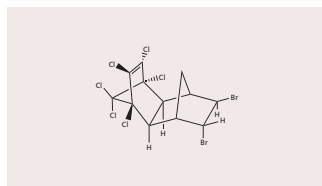
Dec-604



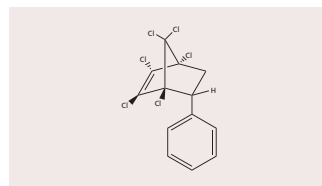
Dec-604CB



CPlus

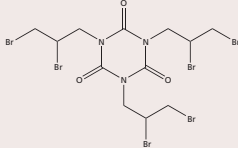


DBALD



HCPN

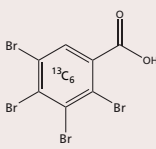
TRIS(2,3-DIBROMOPROPYL)ISOCYANURATE

| Catalogue Number | Product |
|------------------|---|
| T23BPIC |  Tris(2,3-dibromopropyl)isocyanurate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene |

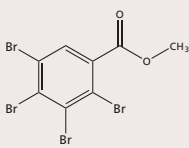
2,3,4,5-TETRABROMOBENZOIC ACID (TBBA)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--------------------------------|-----------------|
| TBBA | 2,3,4,5-Tetrabromobenzoic acid | 1.2 ml 50 µg/ml |

¹³C-LABELLED TETRABROMOBENZOIC ACID (TBBA)

| Catalogue Number | Product |
|------------------|---|
| MTBBA |  2,3,4,5-Tetrabromobenzoic acid [¹³ C ₆ -ring] 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; Isotopic purity; 99% or greater |

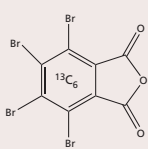
METHYL-2,3,4,5-TETRABROMOBENZOATE (MeTBBA)

| Catalogue Number | Product |
|------------------|---|
| MeTBBA |  Methyl-2,3,4,5-Tetrabromobenzoate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene |

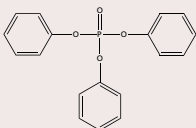
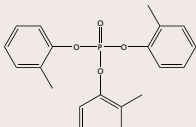
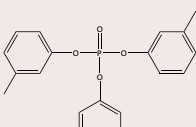
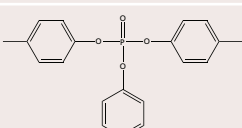
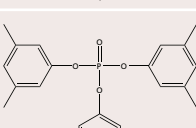
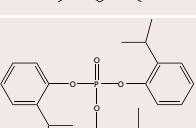
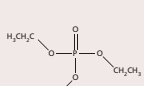
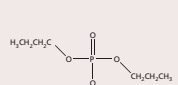
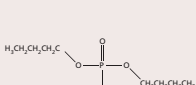
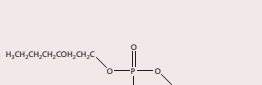
TETRABROMOPHTHALIC ANHYDRIDE (TBPA_n)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|-------------------|------------------------------|-----------------|
| TBPA _n | Tetrabromophthalic anhydride | 1.2 ml 50 µg/ml |

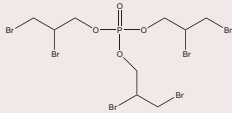
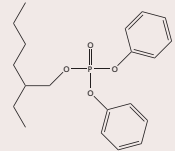
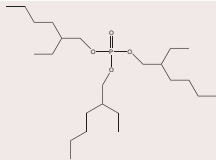
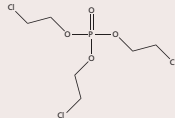
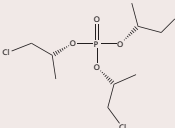
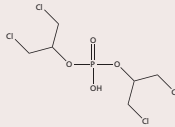
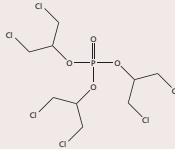
¹³C-LABELLED TETRABROMOPHTHALIC ANHYDRIDE (MTBPA_n)

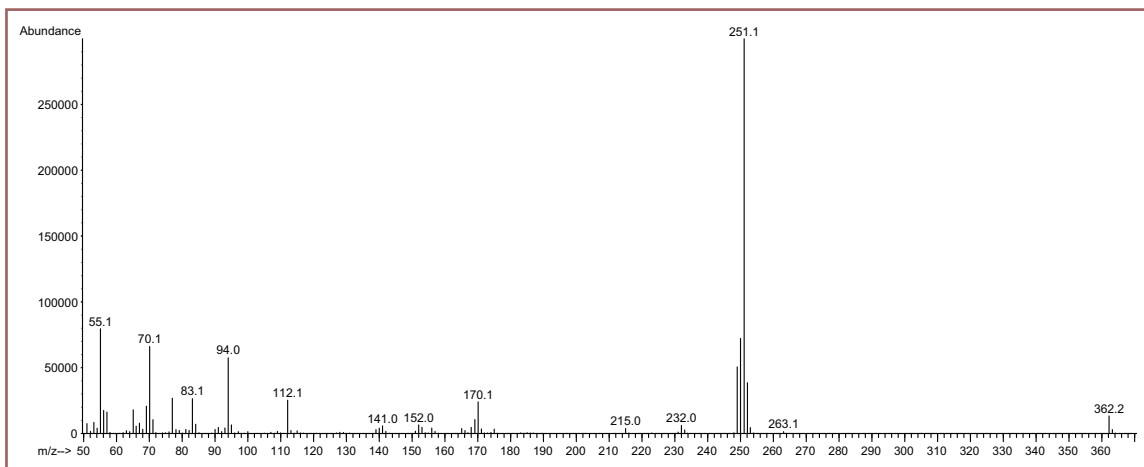
| Catalogue Number | Product |
|--------------------|--|
| MTBPA _n |  Tetrabromo[¹³ C ₆]phthalic anhydride 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater |

NATIVE ORGANOPHOSPHORUS COMPOUNDS

| Catalogue Number | Product |
|------------------|--|
| TPP |  <p>Triphenylphosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TOTP |  <p>Tri-o-tolyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TMTP |  <p>Tri-m-tolyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TPTP |  <p>Tri-p-tolyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| T35DMPP |  <p>Tris(3,5-dimethylphenyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| T2IPPP |  <p>Tris(2-isopropylphenyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TEP |  <p>Tri-ethyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TPrP |  <p>Tri-n-propyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TBP |  <p>Tri-n-butyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TBEP |  <p>Tris(2-butoxyethyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

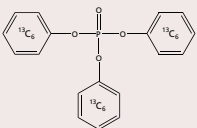
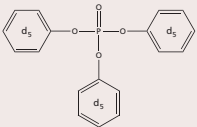
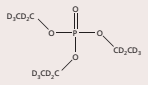
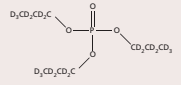
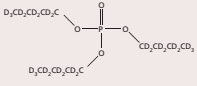
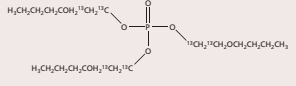
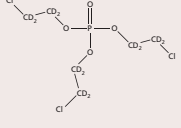
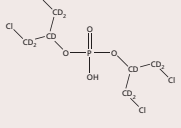
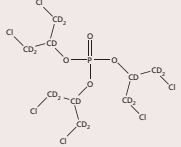
NATIVE ORGANOPHOSPHORUS COMPOUNDS

| Catalogue Number | Product |
|------------------|--|
| TDBPP |  <p>Tris(2,3-dibromopropyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| EHDP |  <p>2-Ethylhexyl diphenyl phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TEHP |  <p>Tris(2-ethylhexyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TCEP |  <p>Tris(2-chloroethyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| TCPP |  <p>Tris[(2<i>R</i>)-1-chloro-2-propyl] phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| BDCP |  <p>Bis(1,3-dichloro-2-propyl) phosphate (97% pure) 1.2 ml; 48.5 µg/ml (±2.4 µg/ml); in <u>acetonitrile</u></p> |
| TDCPP |  <p>Tris(1,3-dichloro-2-propyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |



HRGC/LRMS EI+ Spectra Data for EHDP on a 15m DB-5HT column.

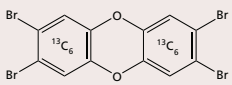
MASS-LABELLED ORGANOPHOSPHORUS COMPOUNDS

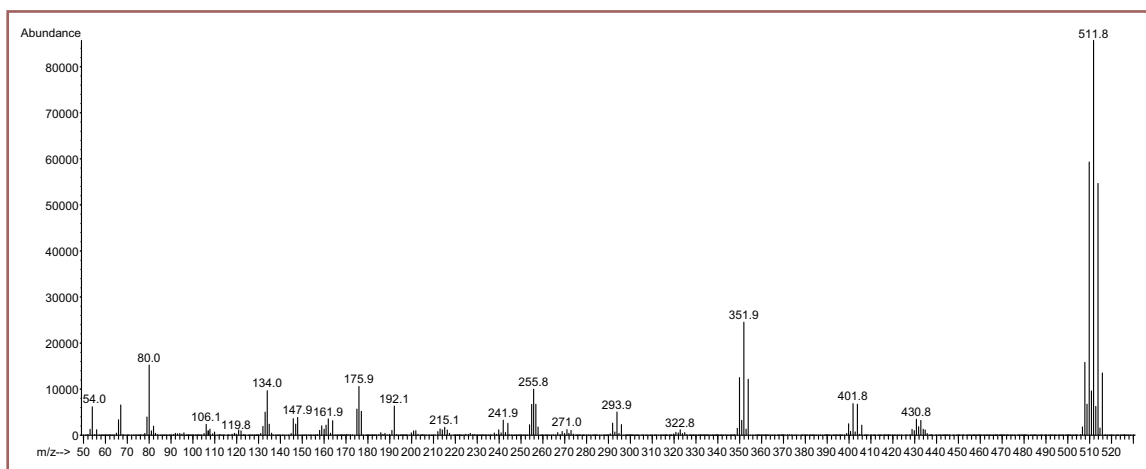
| Catalogue Number | Product |
|------------------|--|
| MTPP |  <p>¹³C₁₈-Triphenylphosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater (¹³C₁₈)</p> |
| dTPP |  <p>Triphenylphosphate-d₁₅ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₁₅)</p> |
| dTEP |  <p>Tri-ethyl phosphate-d₁₅ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₁₅)</p> |
| dTPrP |  <p>Tri-n-propyl phosphate-d₂₁ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₂₁)</p> |
| dTBP |  <p>Tri-n-butyl phosphate-d₂₇ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₂₇)</p> |
| MGTBEP |  <p>Tris(2-butoxy-[¹³C₂]-ethyl) phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene Isotopic purity; 99% or greater (¹³C₆)</p> |
| dTCEP |  <p>Tris(2-chloroethyl) phosphate-d₁₂ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₁₂)</p> |
| dBDCP |  <p>Bis(1,3-dichloro-2-propyl) phosphate-d₁₀ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in <u>acetonitrile</u>; Isotopic purity; 98% or greater (²H₁₀)</p> |
| dTDCPP |  <p>Tris(1,3-dichloro-2-propyl) phosphate-d₁₅ 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 98% or greater (²H₁₅)</p> |

NATIVE BROMINATED DIBENZO-p-DIOXINS (PBDDs)

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|---------------------|--|-----------------|
| BDD-1 | 1-Bromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-27/28 | 2,7/2,8-Dibromodibenzo-p-dioxin mix | 1.2 ml 50 µg/ml |
| BDD-237 | 2,3,7-Tribromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-1234 | 1,2,3,4-Tetrabromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-1378 | 1,3,7,8-Tetrabromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-2378 | 2,3,7,8-Tetrabromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-12378 | 1,2,3,7,8-Pentabromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-12478 | 1,2,4,7,8-Pentabromodibenzo-p-dioxin | 1.2 ml 50 µg/ml |
| BDD-1234678 | 1,2,3,4,6,7,8-Heptabromodibenzo-p-dioxin | 1.2 ml 25 µg/ml |
| BDD-12346789 | Octabromodibenzo-p-dioxin | 1.2 ml 10 µg/ml |

MASS-LABELLED BROMINATED DIBENZO-p-DIOXIN

| Catalogue Number | Product |
|------------------|---|
| MBDD-2378 | <div style="display: flex; align-items: center;"> <div style="margin-right: 20px;">  </div> <div> <p>2,3,7,8-Tetrabromo[¹³C₁₂]dibenzo-p-dioxin</p> <p>1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene;</p> <p>Isotopic purity; 99% or greater</p> </div> </div> |



HRGC/LRMS EI+ Spectra Data for MBDD-2378 on a 15m DB-5HT column.

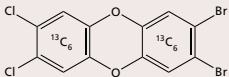
NATIVE BROMINATED DIBENZOFURANS (PBDFs)

| Catalogue Number | Product (toluene solution) | Qty/Conc | |
|---------------------|--------------------------------------|----------|----------|
| BDF-4 | 4-Bromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-24 | 2,4-Dibromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-28 | 2,8-Dibromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-138 | 1,3,8-Tribromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-234 | 2,3,4-Tribromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-238 | 2,3,8-Tribromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-247 | 2,4,7-Tribromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-1278 | 1,2,7,8-Tetrabromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-2378 | 2,3,7,8-Tetrabromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-12378 | 1,2,3,7,8-Pentabromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-23478 | 2,3,4,7,8-Pentabromodibenzofuran | 1.2 ml | 50 µg/ml |
| BDF-1234678 | 1,2,3,4,6,7,8-Heptabromodibenzofuran | 1.2 ml | 25 µg/ml |
| BDF-12346789 | Octabromodibenzofuran (96%) | 1.2 ml | 24 µg/ml |

NATIVE BROMO/CHLORO DIBENZO-p-DIOXINS

| Catalogue Number | Product (toluene solution) | Qty/Conc | |
|---------------------|---|----------|----------|
| 7-B-23-CDD | 7-Bromo-2,3-dichlorodibenzo-p-dioxin (96%) | 1.2 ml | 48 µg/ml |
| 2-B-378-CDD | 2-Bromo-3,7,8-trichlorodibenzo-p-dioxin (96%) | 1.2 ml | 48 µg/ml |
| 2-B-1378-CDD | 2-Bromo-1,3,7,8-tetrachlorodibenzo-p-dioxin (96%) | 1.2 ml | 48 µg/ml |
| 23-B-78-CDD | 2,3-Dibromo-7,8-dichlorodibenzo-p-dioxin | 1.2 ml | 50 µg/ml |

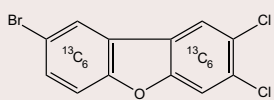
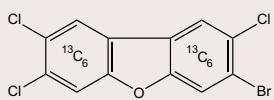
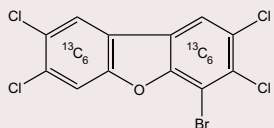
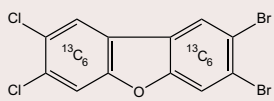
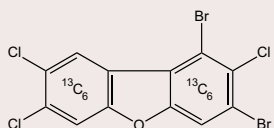
MASS-LABELLED BROMO/CHLORO DIBENZO-p-DIOXIN

| Catalogue Number | Product |
|---------------------|--|
| M23-B-78-CDD | <div style="display: flex; align-items: center;">  <div style="margin-left: 10px;"> <p>2,3-Dibromo-7,8-dichloro[¹³C₁₂]dibenzo-p-dioxin</p> <p>1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene;</p> <p>Isotopic purity; 99% or greater</p> </div> </div> |

NATIVE BROMO/CHLORO DIBENZOFURANS

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|---------------------|---|-----------------|
| 8-B-23-CDF | 8-Bromo-2,3-dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| 3-B-278-CDF | 3-Bromo-2,7,8-trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| 8-B-234-CDF | 8-Bromo-2,3,4-trichlorodibenzofuran | 1.2 ml 50 µg/ml |
| 4-B-2378-CDF | 4-Bromo-2,3,7,8-tetrachlorodibenzofuran | 1.2 ml 50 µg/ml |
| 12-B-78-CDF | 1,2-Dibromo-7,8-dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| 23-B-78-CDF | 2,3-Dibromo-7,8-dichlorodibenzofuran | 1.2 ml 50 µg/ml |
| 13-B-278-CDF | 1,3-Dibromo-2,7,8-trichlorodibenzofuran | 1.2 ml 50 µg/ml |

MASS-LABELLED BROMO/CHLORO DIBENZOFURANS

| Catalogue Number | Product |
|----------------------|--|
| M8-B-23-CDF |  <p>8-Bromo-2,3-dichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |
| M3-B-278-CDF |  <p>3-Bromo-2,7,8-trichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |
| M4-B-2378-CDF |  <p>4-Bromo-2,3,7,8-tetrachloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |
| M23-B-78-CDF |  <p>2,3-Dibromo-7,8-dichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |
| M13-B-278-CDF |  <p>1,3-Dibromo-2,7,8-trichloro[¹³C₁₂]dibenzofuran 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic purity; 99% or greater</p> |

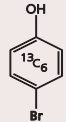
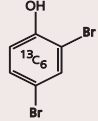
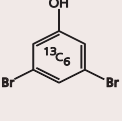
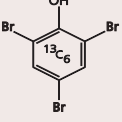
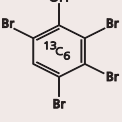
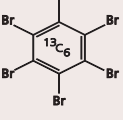
NATIVE BROMOPHENOLS

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|----------------------------|------------------|
| BRP-2 | 2-Bromophenol | 1.2 ml 100 µg/ml |
| BRP-3 | 3-Bromophenol | 1.2 ml 100 µg/ml |
| BRP-4 | 4-Bromophenol | 1.2 ml 100 µg/ml |
| BRP-23 | 2,3-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-24 | 2,4-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-25 | 2,5-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-26 | 2,6-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-34 | 3,4-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-35 | 3,5-Dibromophenol | 1.2 ml 100 µg/ml |
| BRP-234 | 2,3,4-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-235 | 2,3,5-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-236 | 2,3,6-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-245 | 2,4,5-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-246 | 2,4,6-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-345 | 3,4,5-Tribromophenol | 1.2 ml 100 µg/ml |
| BRP-2345 | 2,3,4,5-Tetrabromophenol | 1.2 ml 100 µg/ml |
| BRP-2346 | 2,3,4,6-Tetrabromophenol | 1.2 ml 100 µg/ml |
| BRP-2356 | 2,3,5,6-Tetrabromophenol | 1.2 ml 100 µg/ml |
| BRP-23456 | Pentabromophenol | 1.2 ml 100 µg/ml |

NATIVE BROMOPHENOLS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|---------------------------------------|----------|
| BRPS | Native Bromophenols; Solution/Mixture | 1.2 ml |
| | 2-Bromophenol | 5 µg/ml |
| | 3-Bromophenol | 5 µg/ml |
| | 4-Bromophenol | 5 µg/ml |
| | 2,3-Dibromophenol | 5 µg/ml |
| | 2,4-Dibromophenol | 5 µg/ml |
| | 2,5-Dibromophenol | 5 µg/ml |
| | 2,6-Dibromophenol | 5 µg/ml |
| | 3,4-Dibromophenol | 5 µg/ml |
| | 3,5-Dibromophenol | 5 µg/ml |
| | 2,3,4-Tribromophenol | 5 µg/ml |
| | 2,3,5-Tribromophenol | 5 µg/ml |
| | 2,3,6-Tribromophenol | 5 µg/ml |
| | 2,4,5-Tribromophenol | 5 µg/ml |
| | 2,4,6-Tribromophenol | 5 µg/ml |
| | 3,4,5-Tribromophenol | 5 µg/ml |
| | 2,3,4,5-Tetrabromophenol | 5 µg/ml |
| | 2,3,4,6-Tetrabromophenol | 5 µg/ml |
| | 2,3,5,6-Tetrabromophenol | 5 µg/ml |
| | Pentabromophenol | 5 µg/ml |

MASS-LABELLED BROMOPHENOLS

| Catalogue Number | Product |
|-------------------|---|
| MBRP-4 |  4-Bromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |
| MBRP-24 |  2,4-Dibromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |
| MBRP-35 |  3,5-Dibromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |
| MBRP-246 |  2,4,6-Tribromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |
| MBRP-2346 |  2,3,4,6-Tetrabromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |
| MBRP-23456 |  Pentabromo[¹³ C ₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in toluene |

* Unless stated otherwise, isotopic purities of these compounds are 99% or greater.

MASS-LABELLED BROMOPHENOLS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|----------|
| MBRPS | Mass-Labelled Bromophenols; Solution/Mixture | 1.2 ml |
| | 4-Bromo[¹³ C ₆]phenol | 5 µg/ml |
| | 2,4-Dibromo[¹³ C ₆]phenol | 5 µg/ml |
| | 2,4,6-Tribromo[¹³ C ₆]phenol | 5 µg/ml |
| | 2,3,4,6-Tetrabromo[¹³ C ₆]phenol | 5 µg/ml |
| | Pentabromo[¹³ C ₆]phenol | 5 µg/ml |



*In order to keep up-to-date on new releases from Wellington Laboratories Inc.,
please visit our web-site at www.well-labs.com.*

New Products are announced in the Wellington Reporter.

Lydia Rennie
Cobourg, Ontario, Canada



PERFLUORINATED COMPOUNDS (PFCs)

Wellington started to synthesize perfluorinated compounds in 2004 and, since then, we have regularly added new native and mass-labelled standards to our inventory. In this section you will find individual standards of the following groups of compounds including, in most cases, mass-labelled analogues as well as some useful solution/mixtures:

PFC-C-CVS Calibration Set and Support Solutions

Perfluoroalkanesulfonates (PFASs)

Perfluoroalkylcarboxylic acids (PFCAs)

Perfluorooctanesulfonamides (FOSAs)

Perfluorooctanesulfonamidoethanols (FOSEs)

Perfluorooctanesulfonamidoacetic acids (FOSAAs)

Fluorinated Telomer Alcohols (FTOHs)

Fluorinated Telomer Acids (FTAs)

Unsaturated Fluorinated Telomer Acids (FTUAs)

Perfluoroalkylphosphonic Acids (PFAPAs)

Perfluoroalkylphosphinic Acids (X:XPFPi)

Polyfluorinated Phosphate Esters (PAPs and SAmPAPs)

Fluorinated Telomer Acrylates and Acetates (FTAcrs and FTOAcS)

PFCs are still emerging environmental contaminants and each of the groups of compounds listed above pose unique analytical challenges. In addition, the individual isomers, such as the branched PFOA and PFOS isomers, are being found to have different toxicokinetic and ecokinetic properties. Thus our inventory of PFCs will continue to grow and we would urge you to visit our website for announcements of new products.



PFC-CVS-C

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|-------------------------------|--------------|
| PFC-CVS-C | PFC-CVS-C | 1 kit |
| | Calibration Solutions CS1-CS5 | (5 ampoules) |
| PFC-C-CS1 | CS1 | 200 µl |
| PFC-C-CS2 | CS2 | 200 µl |
| PFC-C-CS3 | CS3 | 200 µl |
| PFC-C-CS4 | CS4 | 200 µl |
| PFC-C-CS5 | CS5 | 200 µl |

| | | PFC-C- CS1 (ng/ml) | PFC-C- CS2 (ng/ml) | PFC-C- CS3 (ng/ml) | PFC-C- CS4 (ng/ml) | PFC-C- CS5 (ng/ml) |
|--|----------|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| Native PFCs | | | | | | |
| Perfluoro-n-butanoic acid | PFBA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-pentanoic acid | PFPeA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-hexanoic acid | PFHxA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-heptanoic acid | PFHpA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-octanoic acid | PFOA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-nonanoic acid | PFNA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-decanoic acid | PFDA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-undecanoic acid | PFUdA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-dodecanoic acid | PFDoA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-tridecanoic acid | PFTrDA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-tetradecanoic acid | PFTeDA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-hexadecanoic acid | PFHxDA | 2.0 | 10 | 50 | 200 | 1000 |
| Perfluoro-n-octadecanoic acid | PFODA | 2.0 | 10 | 50 | 200 | 1000 |
| Potassium perfluoro-1-butanefluorobutanesulfonate | L-PFBS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-pentanesulfonate | L-PFPeS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-hexanesulfonate | L-PFHxS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-heptanesulfonate | L-PFHpS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-octanesulfonate | L-PFOS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-nonanesulfonate | L-PFNS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-decanesulfonate | L-PFDS | 2.0 | 10 | 50 | 200 | 1000 |
| Sodium perfluoro-1-dodecanesulfonate | L-PFDoS | 2.0 | 10 | 50 | 200 | 1000 |
| Mass-Labelled PFCs Extraction Standards | | | | | | |
| Perfluoro-n-[¹³ C ₄]butanoic acid | MPFBA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[¹³ C ₅]pentanoic acid | M5PFPeA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2,3,4,6- ¹³ C ₅]hexanoic acid | M5PFHxA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2,3,4- ¹³ C ₄]heptanoic acid | M4PFHpA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[¹³ C ₈]octanoic acid | M8PFOA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[¹³ C ₉]nonanoic acid | M9PFNA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2,3,4,5,6- ¹³ C ₆]decanoic acid | M6PFDA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2,3,4,5,6,7- ¹³ C ₇]undecanoic acid | M7PFUdA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2- ¹³ C ₂]dodecanoic acid | MPFDoA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2- ¹³ C ₂]tetradecanoic acid | M2PFTeDA | 50 | 50 | 50 | 50 | 50 |
| Sodium perfluoro-1-[2,3,4- ¹³ C ₃]butanesulfonate | M3PFBS | 50 | 50 | 50 | 50 | 50 |
| Sodium perfluoro-1-[1,2,3- ¹³ C ₃]hexanesulfonate | M3PFHxS | 50 | 50 | 50 | 50 | 50 |
| Sodium perfluoro-1-[¹³ C ₈]octanesulfonate | M8PFOS | 50 | 50 | 50 | 50 | 50 |
| Mass-Labelled PFCs Injection Standards | | | | | | |
| Perfluoro-n-[2,3,4- ¹³ C ₃]butanoic acid | M3PFBA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2- ¹³ C ₂]octanoic acid | M2PFOA | 50 | 50 | 50 | 50 | 50 |
| Perfluoro-n-[1,2- ¹³ C ₂]decanoic acid | MPFDA | 50 | 50 | 50 | 50 | 50 |
| Sodium perfluoro-1-[1,2,3,4- ¹³ C ₄]octanesulfonate | MPFOS | 50 | 50 | 50 | 50 | 50 |

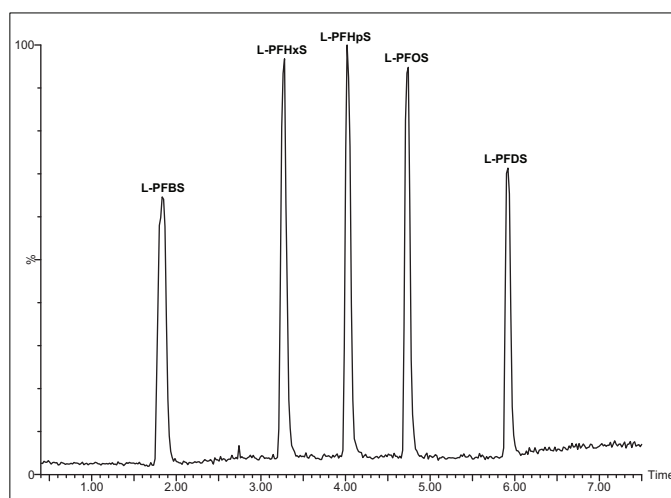
NOTE: All perfluoroalkylsulfonate concentrations are reported as the salt.

NATIVE LINEAR PERFLUOROALKYLSULFONATES (PFASs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---------------------------------------|-----------------|
| L-PFBS | Potassium perfluoro-1-butanesulfonate | 1.2 ml 50 µg/ml |
| L-PFPeS | Sodium perfluoro-1-pentanesulfonate | 1.2 ml 50 µg/ml |
| L-PFHxS | Sodium perfluoro-1-hexanesulfonate | 1.2 ml 50 µg/ml |
| L-PFHpS | Sodium perfluoro-1-heptanesulfonate | 1.2 ml 50 µg/ml |
| L-PFOS | Sodium perfluoro-1-octanesulfonate | 1.2 ml 50 µg/ml |
| L-PFOSK | Potassium perfluoro-1-octanesulfonate | 1.2 ml 50 µg/ml |
| L-PFNS | Sodium perfluoro-1-nonanesulfonate | 1.2 ml 50 µg/ml |
| L-PFDS | Sodium perfluoro-1-decanesulfonate | 1.2 ml 50 µg/ml |
| L-PFDoS | Sodium perfluoro-1-dodecanesulfonate | 1.2 ml 50 µg/ml |

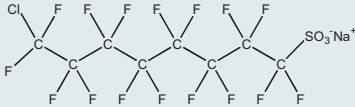
NATIVE PERFLUOROALKYLSULFONATES: SOLUTION/MIXTURE

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---------------------------------------|----------|
| PFS-MXA | Native PFAS Solution/Mixture | 1.2 ml |
| | Potassium perfluoro-1-butanesulfonate | 2 µg/ml |
| | Sodium perfluoro-1-hexanesulfonate | 2 µg/ml |
| | Sodium perfluoro-1-heptanesulfonate | 2 µg/ml |
| | Sodium perfluoro-1-octanesulfonate | 2 µg/ml |
| | Sodium perfluoro-1-decanesulfonate | 2 µg/ml |



Chromatogram of PFS-MXA (Acquity UPLC BEH Shield RP₁₈ column)

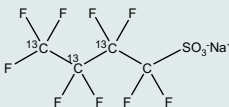
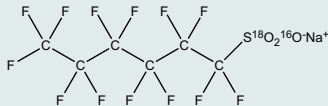
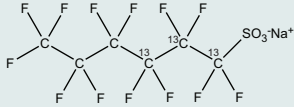
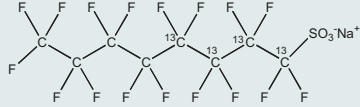
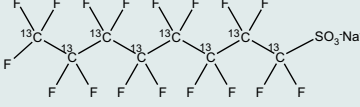
NATIVE CHLORO-PERFLUOROALKYLSULFONATE

| Catalogue Number | Product |
|------------------|--|
| 8Cl-PFOS |  <p>Sodium 8-chloroperfluoro-1-octanesulfonate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |

NATIVE BRANCHED PERFLUOROALKYLSULFONATES

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| br-PFHxSK | L-PFHxS with branched isomers (Potassium Salt) | 1.2 ml 50 µg/ml |
| br-PFOSK | L-PFOSK with branched isomers | 1.2 ml 50 µg/ml |
| T-PFOS | Potassium perfluorooctanesulfonate (Technical Grade) | 1.2 ml 50 µg/ml |
| NaP3MHpS | Sodium perfluoro-3-methylheptanesulfonate | 1.2 ml 50 µg/ml |
| NaP6MHpS | Sodium perfluoro-6-methylheptanesulfonate | 1.2 ml 50 µg/ml |
| ipPFNS | Sodium perfluoro-7-methyloctanesulfonate | 1.2 ml 50 µg/ml |

MASS-LABELLED PERFLUOROALKYLSULFONATES

| Catalogue Number | Product |
|------------------|--|
| M3PFBS |  <p>Sodium perfluoro-1-[2,3,4-¹³C₃]-butanesulfonate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 2,3,4-¹³C₃</p> |
| MPFHxS |  <p>Sodium perfluoro-1-hexane[¹⁸O₂]sulfonate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 94% ¹⁸O₂</p> |
| M3PFHxS |  <p>Sodium perfluoro-1-[1,2,3-¹³C₃]-hexanesulfonate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3-¹³C₃</p> |
| MPFOS |  <p>Sodium perfluoro-1-[1,2,3,4-¹³C₄]-octanesulfonate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4-¹³C₄</p> |
| M8PFOS |  <p>Sodium perfluoro-1-[¹³C₈]-octanesulfonate 1.2 ml; 48.5 µg/ml (±2.4 µg/ml); in methanol >99% linear; >99% ¹³C₈</p> |

NOTE: All perfluoroalkylsulfonate concentrations are reported as the salt.

NATIVE LINEAR PERFLUOROALKYLCARBOXYLIC ACIDS (PFCAs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--------------------------------|-----------------|
| PFBA | Perfluoro-n-butanoic acid | 1.2 ml 50 µg/ml |
| PFPeA | Perfluoro-n-pentanoic acid | 1.2 ml 50 µg/ml |
| PFHxA | Perfluoro-n-hexanoic acid | 1.2 ml 50 µg/ml |
| PFHpA | Perfluoro-n-heptanoic acid | 1.2 ml 50 µg/ml |
| PFOA | Perfluoro-n-octanoic acid | 1.2 ml 50 µg/ml |
| PFNA | Perfluoro-n-nonanoic acid | 1.2 ml 50 µg/ml |
| PFDA | Perfluoro-n-decanoic acid | 1.2 ml 50 µg/ml |
| PFUdA | Perfluoro-n-undecanoic acid | 1.2 ml 50 µg/ml |
| PFDoA | Perfluoro-n-dodecanoic acid | 1.2 ml 50 µg/ml |
| PFTTrDA | Perfluoro-n-tridecanoic acid | 1.2 ml 50 µg/ml |
| PFTeDA | Perfluoro-n-tetradecanoic acid | 1.2 ml 50 µg/ml |
| PFHxDA | Perfluoro-n-hexadecanoic acid | 1.2 ml 50 µg/ml |
| PFODA | Perfluoro-n-octadecanoic acid | 1.2 ml 50 µg/ml |

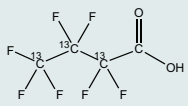
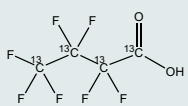
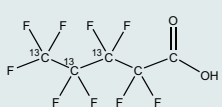

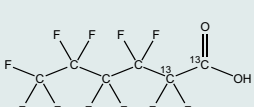
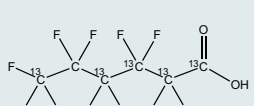
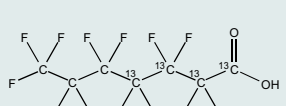
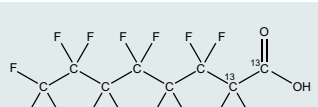
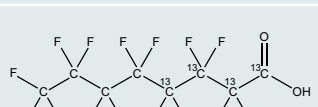
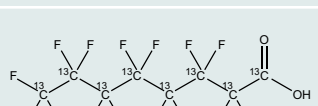
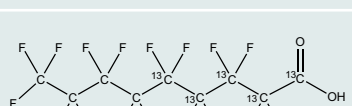
NATIVE PERFLUOROALKYLCARBOXYLIC ACIDS: SOLUTION/MIXTURE

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--------------------------------|----------|
| PFC-MXA | Native PFCa Solution/Mixture | 1.2 ml |
| | Perfluoro-n-butanoic acid | 2 µg/ml |
| | Perfluoro-n-pentanoic acid | 2 µg/ml |
| | Perfluoro-n-hexanoic acid | 2 µg/ml |
| | Perfluoro-n-heptanoic acid | 2 µg/ml |
| | Perfluoro-n-octanoic acid | 2 µg/ml |
| | Perfluoro-n-nonanoic acid | 2 µg/ml |
| | Perfluoro-n-decanoic acid | 2 µg/ml |
| | Perfluoro-n-undecanoic acid | 2 µg/ml |
| | Perfluoro-n-dodecanoic acid | 2 µg/ml |
| | Perfluoro-n-tridecanoic acid | 2 µg/ml |
| | Perfluoro-n-tetradecanoic acid | 2 µg/ml |

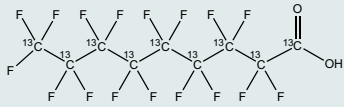
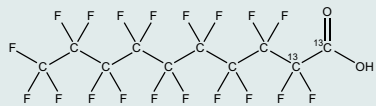
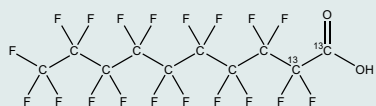
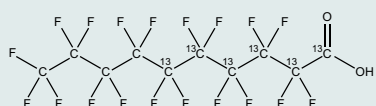
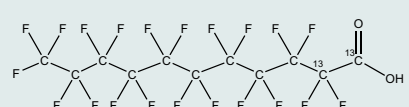
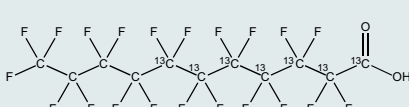
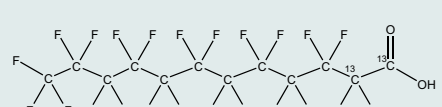
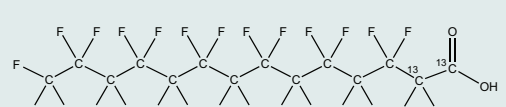

NATIVE BRANCHED PERFLUOROALKYLCARBOXYLIC ACIDS

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| T-PFOA | Perfluorooctanoic acid (Technical Grade) | 1.2 ml 50 µg/ml |
| P3MHpA | Perfluoro-3-methylheptanoic acid | 1.2 ml 50 µg/ml |
| P4MOA | Perfluoro-4-methyloctanoic acid | 1.2 ml 50 µg/ml |
| ipPFNA | Perfluoro-7-methyloctanoic acid (90%) | 1.2 ml 45 µg/ml |
| P355TMHxA | Perfluoro-3,5,5-trimethylhexanoic acid | 1.2 ml 50 µg/ml |
| P37DMOA | Perfluoro-3,7-dimethyloctanoic acid | 1.2 ml 50 µg/ml |

MASS-LABELLED PERFLUOROALKYLCARBOXYLIC ACIDS

| Catalogue Number | Product |
|------------------|--|
| M3PFBA |  <p>Perfluoro-n-[2,3,4-¹³C₃]butanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 2,3,4-¹³C₃</p> |
| MPFBA |  <p>Perfluoro-n-[¹³C₄]butanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% ¹³C₄</p> |
| M3PFPeA |  <p>Perfluoro-n-[3,4,5-¹³C₃]pentanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 3,4,5-¹³C₃</p> |
| M5PFPeA |  <p>Perfluoro-n-[¹³C₅]pentanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% ¹³C₅</p> |
| MPFHxA |  <p>Perfluoro-n-[1,2-¹³C₂]hexanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2-¹³C₂</p> |
| M5PFHxA |  <p>Perfluoro-n-[1,2,3,4,6-¹³C₅]hexanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4,6-¹³C₅</p> |
| M4PFHpA |  <p>Perfluoro-n-[1,2,3,4-¹³C₄]heptanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4-¹³C₄</p> |
| M2PFOA |  <p>Perfluoro-n-[1,2-¹³C₂]octanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2-¹³C₂</p> |
| MPFOA |  <p>Perfluoro-n-[1,2,3,4-¹³C₄]octanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4-¹³C₄</p> |
| M8PFOA |  <p>Perfluoro-n-[¹³C₈]octanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 97.9% ¹³C₈ and 2.1% ¹³C₄</p> |
| MPFNA |  <p>Perfluoro-n-[1,2,3,4,5-¹³C₅]nonanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4,5-¹³C₅</p> |

MASS-LABELLED PERFLUOROALKYLCARBOXYLIC ACIDS

| Catalogue Number | Product |
|--|--|
| M9PFNA  | Perfluoro-n-[¹³ C ₉]nonanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% ¹³ C ₉ |
| MPFDA  | Perfluoro-n-[1,2- ¹³ C ₂]decanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2- ¹³ C ₂ |
| MPFDA-A  | Perfluoro-n-[1,2- ¹³ C ₂]decanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in <u>acetonitrile</u> >99% linear; >99% 1,2- ¹³ C ₂ |
| M6PFDA  | Perfluoro-n-[1,2,3,4,5,6- ¹³ C ₆]decanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4,5,6- ¹³ C ₆ |
| MPFUdA  | Perfluoro-n-[1,2- ¹³ C ₂]undecanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2- ¹³ C ₂ |
| M7PFUdA  | Perfluoro-n-[1,2,3,4,5,6,7- ¹³ C ₇]undecanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2,3,4,5,6,7- ¹³ C ₇ |
| MPFDoA  | Perfluoro-n-[1,2- ¹³ C ₂]dodecanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2- ¹³ C ₂ |
| M2PFTeDA  | Perfluoro-n-[1,2- ¹³ C ₂]tetradecanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2- ¹³ C ₂ |
| M2PFHxDA  | Perfluoro-n-[1,2- ¹³ C ₂]hexadecanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; >99% 1,2- ¹³ C ₂ |

MIXED NATIVE PFCAs AND PFASs: SOLUTION/MIXTURES

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|---|---|----------|
| PFAC-MXA | Native PFCAs and PFASs Solution/Mixture | 1.2 ml |
| Perfluoro-n-butanoic acid | | 5 µg/ml |
| Perfluoro-n-pentanoic acid | | 5 µg/ml |
| Perfluoro-n-hexanoic acid | | 5 µg/ml |
| Perfluoro-n-heptanoic acid | | 5 µg/ml |
| Perfluoro-n-octanoic acid | | 5 µg/ml |
| Perfluoro-n-nonanoic acid | | 5 µg/ml |
| Perfluoro-n-decanoic acid | | 5 µg/ml |
| Potassium perfluoro-1-butanedisulfonate | | 5 µg/ml |
| Sodium perfluoro-1-hexanesulfonate | | 5 µg/ml |
| Sodium perfluoro-1-octanesulfonate | | 5 µg/ml |
| PFAC-MXB | Native PFCAs and PFASs Solution/Mixture | 1.2 ml |
| Perfluoro-n-butanoic acid | | 2 µg/ml |
| Perfluoro-n-pentanoic acid | | 2 µg/ml |
| Perfluoro-n-hexanoic acid | | 2 µg/ml |
| Perfluoro-n-heptanoic acid | | 2 µg/ml |
| Perfluoro-n-octanoic acid | | 2 µg/ml |
| Perfluoro-n-nonanoic acid | | 2 µg/ml |
| Perfluoro-n-decanoic acid | | 2 µg/ml |
| Perfluoro-n-undecanoic acid | | 2 µg/ml |
| Perfluoro-n-dodecanoic acid | | 2 µg/ml |
| Perfluoro-n-tridecanoic acid | | 2 µg/ml |
| Perfluoro-n-tetradecanoic acid | | 2 µg/ml |
| Perfluoro-n-hexadecanoic acid | | 2 µg/ml |
| Perfluoro-n-octadecanoic acid | | 2 µg/ml |
| Potassium perfluoro-1-butanedisulfonate | | 2 µg/ml |
| Sodium perfluoro-1-hexanesulfonate | | 2 µg/ml |
| Sodium perfluoro-1-octanesulfonate | | 2 µg/ml |
| Sodium perfluoro-1-decanedisulfonate | | 2 µg/ml |

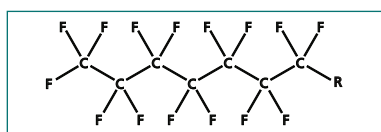
MIXED MASS-LABELLED PFCAs AND PFASs: SOLUTION/MIXTURE

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|--|--|----------|
| MPFAC-MXA | Mass-Labelled PFCAs and PFASs Solution/Mixture | 1.2 ml |
| Perfluoro-n-[1,2,3,4- ¹³ C ₄]butanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2- ¹³ C ₂]hexanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2,3,4- ¹³ C ₄]octanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2,3,4,5- ¹³ C ₅]nonanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2- ¹³ C ₂]decanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2- ¹³ C ₂]undecanoic acid | | 2 µg/ml |
| Perfluoro-n-[1,2- ¹³ C ₂]dodecanoic acid | | 2 µg/ml |
| Sodium perfluoro-1-hexane[¹⁸ O ₂]sulfonate | | 2 µg/ml |
| Sodium perfluoro-1-[1,2,3,4- ¹³ C ₄]octanesulfonate | | 2 µg/ml |

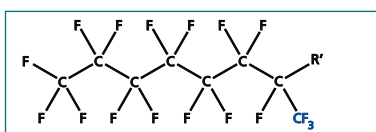
NOTE: Listed concentrations for the perfluoroalkylsulfonates are reported as the salt.

PFOS/PFOA ISOMERS

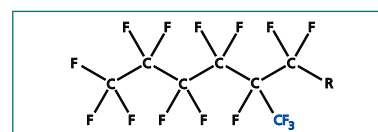
| Catalogue Number | Product (methanol solution) | Qty/Anion Conc |
|------------------|---|-------------------|
| P1MHpS | Sodium Perfluoro-1-methylheptane sulfonate | 200 µl 1.00 µg/ml |
| P3MHpS | Sodium Perfluoro-3-methylheptane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-3-methylheptanoic acid | 1.90 µg/ml |
| P4MHpS | Sodium Perfluoro-4-methylheptane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-4-methylheptanoic acid | 2.20 µg/ml |
| P5MHpS | Sodium Perfluoro-5-methylheptane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-5-methylheptanoic acid | 1.96 µg/ml |
| P6MHpS | Sodium Perfluoro-6-methylheptane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-6-methylheptanoic acid | 3.10 µg/ml |
| P55DMHxS | Sodium Perfluoro-5,5-dimethylhexane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-5,5-dimethylhexanoic acid | 1.95 µg/ml |
| P44DMHxS | Sodium Perfluoro-4,4-dimethylhexane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-4,4-dimethylhexanoic acid | 3.14 µg/ml |
| P45DMHxS | Sodium Perfluoro-4,5-dimethylhexane sulfonate | 200 µl 1.00 µg/ml |
| | Perfluoro-4,5-dimethylhexanoic acid | 1.22 µg/ml |
| | Sodium Perfluoro-3,5-dimethylhexane sulfonate | 0.50 µg/ml |
| | Perfluoro-3,5-dimethylhexanoic acid | 0.60 µg/ml |



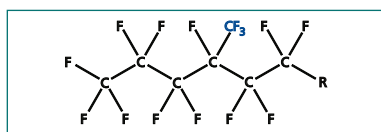
Linear



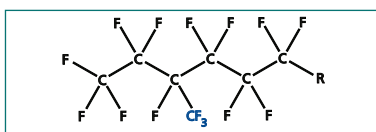
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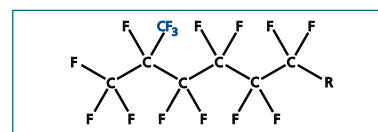
Perfluoro-3-methyl-



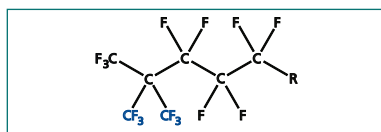
Perfluoro-4-methyl-



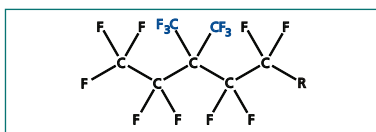
Perfluoro-5-methyl-



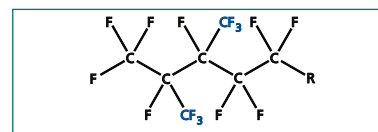
Perfluoro-6-methyl-



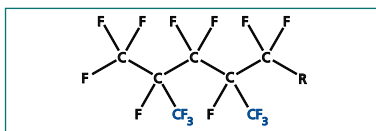
Perfluoro-5,5-dimethyl-



Perfluoro-4,4-dimethyl-



Perfluoro-4,5-dimethyl-



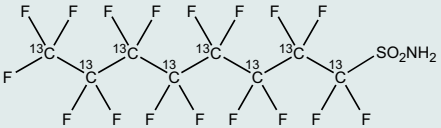
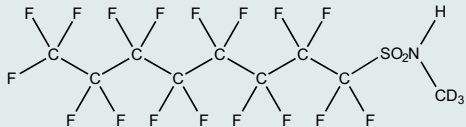
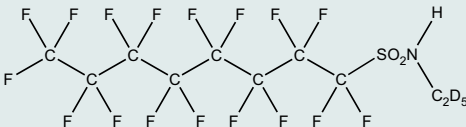
Perfluoro-3,5-dimethyl-

NOTE: R= CO₂⁻ and CF₂SO₃⁻
R'= SO₃⁻ only

NATIVE PERFLUOROCTANESULFONAMIDES (FOSAs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|----------------------|--|-----------------|
| FOSA-I | Perfluoro-1-octanesulfonamide (in isopropanol) | 1.2 ml 50 µg/ml |
| N-MeFOSA-M | N-methylperfluoro-1-octanesulfonamide | 1.2 ml 50 µg/ml |
| N,N-Me2FOSA-M | N,N-dimethylperfluoro-1-octanesulfonamide | 1.2 ml 50 µg/ml |
| N-EtFOSA-M | N-ethylperfluoro-1-octanesulfonamide | 1.2 ml 50 µg/ml |

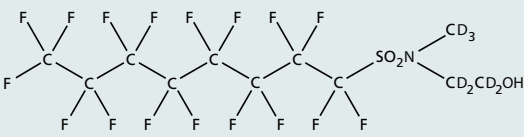
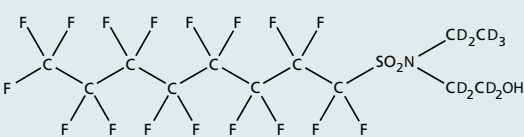
MASS-LABELLED PERFLUOROCTANESULFONAMIDES

| Catalogue Number | Product |
|---------------------|---|
| M8FOSA-I |  <p>Perfluoro-1-[¹³C₈]octanesulfonamide 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol >99% linear; >99% ¹³C₈</p> |
| d-N-MeFOSA-M |  <p>N-methyl-d₃-perfluoro-1-octanesulfonamide 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₃</p> |
| d-N-EtFOSA-M |  <p>N-ethyl-d₅-perfluoro-1-octanesulfonamide 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₅</p> |

NATIVE PERFLUOROCTANESULFONAMIDOETHANOLS (FOSEs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|-------------------|---|-----------------|
| N-MeFOSE-M | 2-(N-methylperfluoro-1-octanesulfonamido)-ethanol | 1.2 ml 50 µg/ml |
| N-EtFOSE-M | 2-(N-ethylperfluoro-1-octanesulfonamido)-ethanol | 1.2 ml 50 µg/ml |

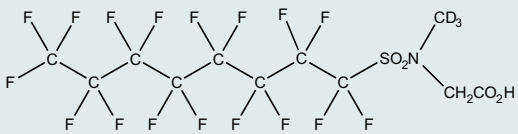
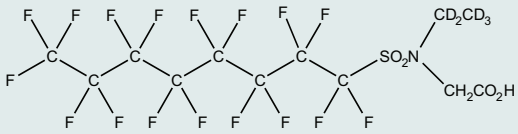
MASS-LABELLED PERFLUOROCTANESULFONAMIDOETHANOLS

| Catalogue Number | Product |
|----------------------|--|
| d7-N-MeFOSE-M |  <p>2-(N-methyl-d₃-perfluoro-1-octanesulfonamido)ethan-d₄-ol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₇</p> |
| d9-N-EtFOSE-M |  <p>2-(N-ethyl-d₅-perfluoro-1-octanesulfonamido)ethan-d₄-ol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₉</p> |

NATIVE PERFLUOROOCETANESULFONAMIDOACETIC ACIDS (FOSAAs)

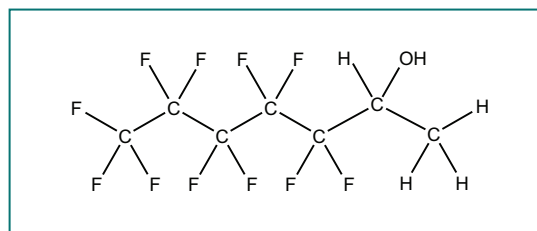
| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---|-----------------|
| FOSAA | Perfluoro-1-octanesulfonamidoacetic acid (in isopropanol) | 1.2 ml 50 µg/ml |
| N-MeFOSAA | N-methylperfluoro-1-octanesulfonamidoacetic acid | 1.2 ml 50 µg/ml |
| N-EtFOSAA | N-ethylperfluoro-1-octanesulfonamidoacetic acid | 1.2 ml 50 µg/ml |

MASS-LABELLED PERFLUOROOCETANESULFONAMIDOACETIC ACIDS

| Catalogue Number | Product |
|---------------------|---|
| d3-N-MeFOSAA |  <p>N-methyl-d₃-perfluoro-1-octanesulfonamidoacetic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₃</p> |
| d5-N-EtFOSAA |  <p>N-ethyl-d₅-perfluoro-1-octanesulfonamidoacetic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol >99% linear; 98% ²H₅</p> |

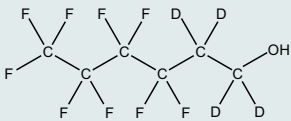
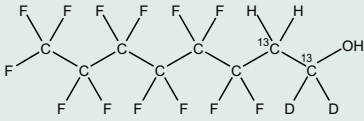
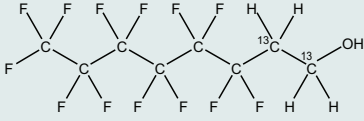
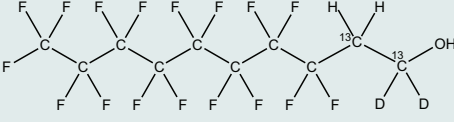
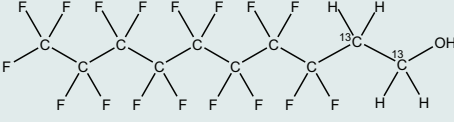
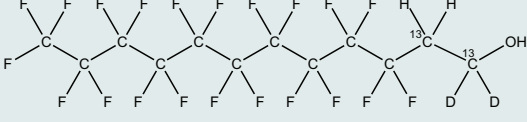
NATIVE TELOMER ALCOHOLS (FTOHs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| FBET | 2-Perfluorobutyl ethanol (4:2) | 1.2 ml 50 µg/ml |
| 5:2sFTOH | 1-Perfluoropentyl ethanol (5:2 secondary) | 1.2 ml 50 µg/ml |
| FHET | 2-Perfluorohexyl ethanol (6:2) | 1.2 ml 50 µg/ml |
| 7:2sFTOH | 1-Perfluoroheptyl ethanol (7:2 secondary) | 1.2 ml 50 µg/ml |
| FOET | 2-Perfluorooctyl ethanol (8:2) | 1.2 ml 50 µg/ml |
| FDET | 2-Perfluorodecyl ethanol (10:2) | 1.2 ml 50 µg/ml |



5:2sFTOH: 1-Perfluoropentyl ethanol

MASS-LABELLED TELOMER ALCOHOLS

| Catalogue Number | Product |
|------------------|---|
| MF BET |  <p>2-Perfluorobutyl-[1,1,2,2-²H₄]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 98% 1,1,2,2-²H₄</p> |
| MF HET |  <p>2-Perfluorohexyl-[1,1-²H₂]-[1,2-¹³C₂]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; 98% 1,1-²H₂, >99% 1,2-¹³C₂</p> |
| M2 FHET |  <p>2-Perfluorohexyl-[1,2-¹³C₂]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; >99% 1,2-¹³C₂</p> |
| MFO ET |  <p>2-Perfluorooctyl-[1,1-²H₂]-[1,2-¹³C₂]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; 98% 1,1-²H₂, >99% 1,2-¹³C₂</p> |
| M2 FO ET |  <p>2-Perfluorooctyl-[1,2-¹³C₂]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; >99% 1,2-¹³C₂</p> |
| MF DE T |  <p>2-Perfluorodecyl-[1,1-²H₂]-[1,2-¹³C₂]-ethanol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; 98% 1,1-²H₂, >99% 1,2-¹³C₂</p> |

NATIVE TELOMER ACIDS (FTAs)

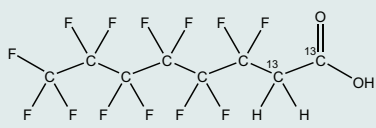
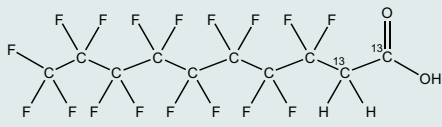
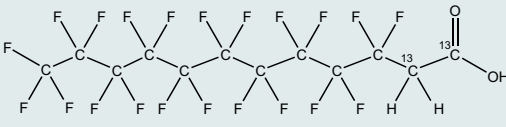
| Catalogue Number | Product (isopropanol solution) | Qty/Conc |
|------------------|--|-----------------|
| FHEA | 2-Perfluorohexyl ethanoic acid (6:2) | 1.2 ml 50 µg/ml |
| FOEA | 2-Perfluorooctyl ethanoic acid (8:2) | 1.2 ml 50 µg/ml |
| FDEA | 2-Perfluorodecyl ethanoic acid (10:2) | 1.2 ml 50 µg/ml |

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---|-----------------|
| FPrPA | 3-Perfluoropropyl propanoic acid (3:3) | 1.2 ml 50 µg/ml |
| FPePA | 3-Perfluoropentyl propanoic acid (5:3) | 1.2 ml 50 µg/ml |
| FHpPA | 3-Perfluoroheptyl propanoic acid (7:3) | 1.2 ml 50 µg/ml |

NATIVE TELOMER ACIDS: SOLUTION/MIXTURE

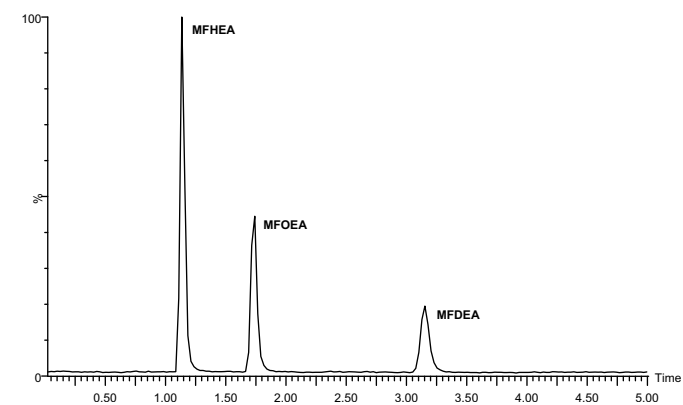
| Catalogue Number | Product (isopropanol solution) | Qty/Conc |
|------------------|--|----------|
| FTA-MXA | Native FTA Solution/Mixture | 1.2 ml |
| | 2-Perfluorohexyl ethanoic acid (6:2) | 2 µg/ml |
| | 2-Perfluorooctyl ethanoic acid (8:2) | 2 µg/ml |
| | 2-Perfluorodecyl ethanoic acid (10:2) | 2 µg/ml |

MASS-LABELLED TELOMER ACIDS

| Catalogue Number | Product |
|------------------|--|
| MFHEA |  <p>2-Perfluorohexyl-[1,2-¹³C₂]-ethanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |
| MFOEA |  <p>2-Perfluorooctyl-[1,2-¹³C₂]-ethanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |
| MFDEA |  <p>2-Perfluorodecyl-[1,2-¹³C₂]-ethanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |

MASS-LABELLED TELOMER ACIDS: SOLUTION/MIXTURE

| Catalogue Number | Product (isopropanol solution) | Qty/Conc |
|------------------|---|----------|
| MFTA-MXA | Mass-Labelled FTA Solution/Mixture | 1.2 ml |
| | 2-Perfluorohexyl-[1,2- ¹³ C ₂]-ethanoic acid (6:2) | 2 µg/ml |
| | 2-Perfluorooctyl-[1,2- ¹³ C ₂]-ethanoic acid (8:2) | 2 µg/ml |
| | 2-Perfluorodecyl-[1,2- ¹³ C ₂]-ethanoic acid (10:2) | 2 µg/ml |

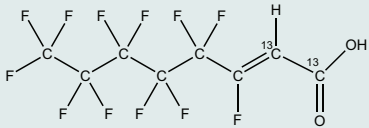
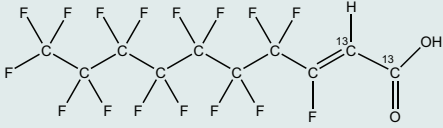
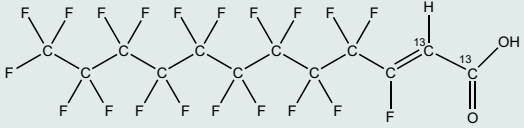


Chromatogram of MFTA-MXA (Acquity UPLC BEH Shield RP₁₈ column)

NATIVE UNSATURATED TELOMER ACIDS (FTUAs)

| Catalogue Number | Product (isopropanol solution) | Qty/Conc |
|------------------|--|-----------------|
| FHUEA | 2H-Perfluoro-2-octenoic acid (6:2) | 1.2 ml 50 µg/ml |
| FOUEA | 2H-Perfluoro-2-decenoic acid (8:2) | 1.2 ml 50 µg/ml |
| FDUEA | 2H-Perfluoro-2-dodecenoic acid (10:2) | 1.2 ml 50 µg/ml |

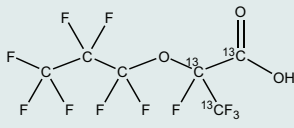
MASS-LABELLED UNSATURATED TELOMER ACIDS

| Catalogue Number | Product |
|------------------|--|
| MFHUEA |  <p>2H-Perfluoro-[1,2-¹³C₂]-2-octenoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |
| MFOUEA |  <p>2H-Perfluoro-[1,2-¹³C₂]-2-decenoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |
| MFDEUA |  <p>2H-Perfluoro-[1,2-¹³C₂]-2-dodecenoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in isopropanol; > 99% ¹³C₂</p> |

NATIVE HEXAFLUOROPROPYLENE OXIDE DIMER ACID

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| HFPO-DA | 2,3,3,3-Tetrafluoro-2-(1,1,2,2,3,3,3-heptafluoropropoxy)propanoic acid | 1.2 ml 50 µg/ml |

MASS-LABELLED HEXAFLUOROPROPYLENE OXIDE DIMER ACID

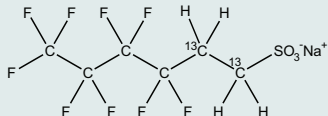
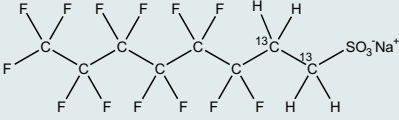
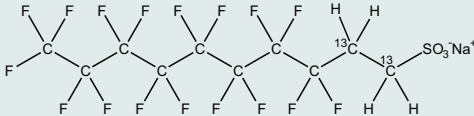
| Catalogue Number | Product |
|------------------|---|
| M3HFPO-DA |  <p>2,3,3,3-Tetrafluoro-2-(1,1,2,2,3,3,3-heptafluoropropoxy)-¹³C₃-propanoic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₃</p> |

NATIVE TELOMER SULFONATES (FTSs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| 4:2FTS | Sodium 1H,1H,2H,2H-perfluorohexane sulfonate (4:2) | 1.2 ml 50 µg/ml |
| 6:2FTS | Sodium 1H,1H,2H,2H-perfluorooctane sulfonate (6:2) | 1.2 ml 50 µg/ml |
| 8:2FTS | Sodium 1H,1H,2H,2H-perfluorodecane sulfonate (8:2) | 1.2 ml 50 µg/ml |
| 10:2FTS | Sodium 1H,1H,2H,2H-perfluorododecane sulfonate (10:2) | 1.2 ml 50 µg/ml |

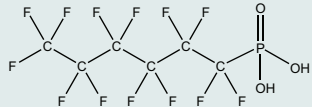
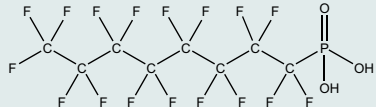
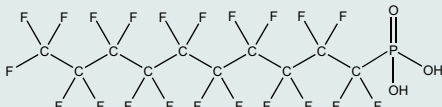
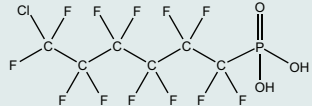
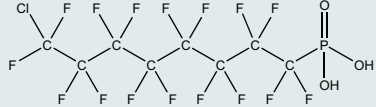
NOTE: Listed concentrations are reported as the salt.

MASS-LABELLED TELOMER SULFONATES

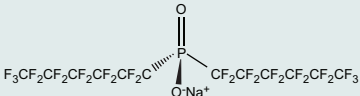
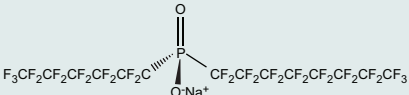
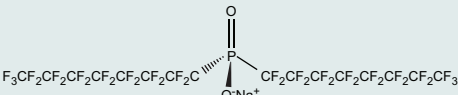
| Catalogue Number | Product |
|------------------|--|
| M2-4:2FTS |  <p>Sodium 1H,1H,2H,2H-perfluoro-1-[1,2-¹³C₂]-hexane sulfonate (4:2) 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₂</p> |
| M2-6:2FTS |  <p>Sodium 1H,1H,2H,2H-perfluoro-1-[1,2-¹³C₂]-octane sulfonate (6:2) 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₂</p> |
| M2-8:2FTS |  <p>Sodium 1H,1H,2H,2H-perfluoro-1-[1,2-¹³C₂]-decane sulfonate (8:2) 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₂</p> |

NOTE: Listed concentration is reported as the salt.

NATIVE PERFLUOROALKYLPHOSPHONIC ACIDS (PFAPAs)

| Catalogue Number | Product |
|------------------|--|
| PFHxPA |  <p>Perfluorohexylphosphonic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| PFOPA |  <p>Perfluorooctylphosphonic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| PFDPA |  <p>Perfluorodecylphosphonic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| Cl-PFHxPA |  <p>6-Chloroperfluorohexylphosphonic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| Cl-PFOPA |  <p>8-Chloroperfluorooctylphosphonic acid 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |

NATIVE SODIUM PERFLUOROALKYL PHOSPHINATES (X:XPFPi)

| Catalogue Number | Product |
|------------------|---|
| 6:6PFPi |  <p>Sodium bis(perfluorohexyl)phosphinate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| 6:8PFPi |  <p>Sodium perfluorohexylperfluorooctylphosphinate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |
| 8:8PFPi |  <p>Sodium bis(perfluorooctyl)phosphinate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol;</p> |

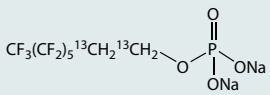
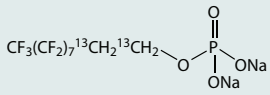
NATIVE MONO-SUBSTITUTED POLYFLUORINATED PHOSPHATE ESTERS

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| 6:2PAP | Sodium 1H,1H,2H,2H-perfluorooctylphosphate | 1.2 ml 50 µg/ml |
| 8:2PAP | Sodium 1H,1H,2H,2H-perfluorodecylphosphate | 1.2 ml 50 µg/ml |

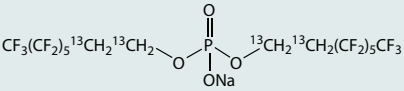
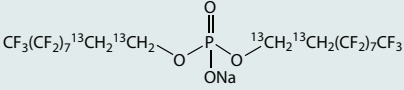
NATIVE DI-SUBSTITUTED POLYFLUORINATED PHOSPHATE ESTERS

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|---------------------|---|-----------------|
| 6:2diPAP | Sodium bis(1H,1H,2H,2H-perfluorooctyl)phosphate | 1.2 ml 50 µg/ml |
| 6:2/8:2diPAP | Sodium (1H,1H,2H,2H-perfluorooctyl-1H,1H,2H,2H-perfluorodecyl)phosphate | 1.2 ml 50 µg/ml |
| 8:2diPAP | Sodium bis(1H,1H,2H,2H-perfluorodecyl)phosphate | 1.2 ml 50 µg/ml |

MASS-LABELLED MONO-SUBSTITUTED POLYFLUORINATED PHOSPHATE ESTERS (PAPs)

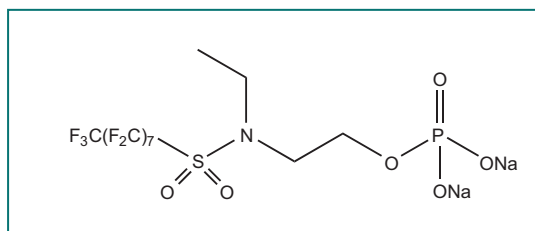
| Catalogue Number | Product |
|------------------|---|
| M2-6:2PAP |  <p>Sodium 1H,1H,2H,2H-[1,2-¹³C₂]perfluorooctylphosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₂</p> |
| M2-8:2PAP |  <p>Sodium 1H,1H,2H,2H-[1,2-¹³C₂]perfluorodecylphosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³C₂</p> |

MASS-LABELLED DI-SUBSTITUTED POLYFLUORINATED PHOSPHATE ESTERS (diPAPs)

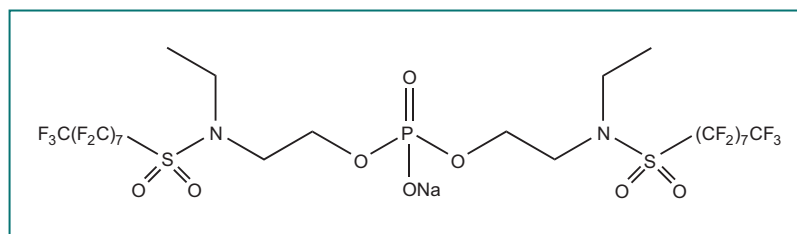
| Catalogue Number | Product |
|---|---|
| M4-6:2diPAP  | Sodium bis(1H,1H,2H,2H-[1,2- ¹³ C ₂]perfluorooctyl)- phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³ C ₄ |
| M4-8:2diPAP  | Sodium bis(1H,1H,2H,2H-[1,2- ¹³ C ₂]perfluorodecyl)- phosphate 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; > 99% ¹³ C ₄ |

NATIVE POLYFLUORINATED PHOSPHATE ESTERS (SAmPAPs)

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---|-----------------|
| SAmPAP | Sodium 2-(N-ethylperfluorooctane-1-sulfonamido)ethyl phosphate | 1.2 ml 50 µg/ml |
| diSAmPAP | Sodium bis[2-(N-ethylperfluorooctane-1-sulfonamido)ethyl] phosphate | 1.2 ml 50 µg/ml |



SAmPAP



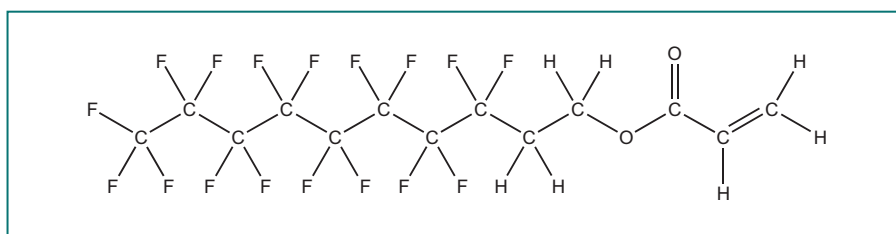
diSAmPAP

NATIVE TELOMER ACRYLATES (X:2FTAcr)

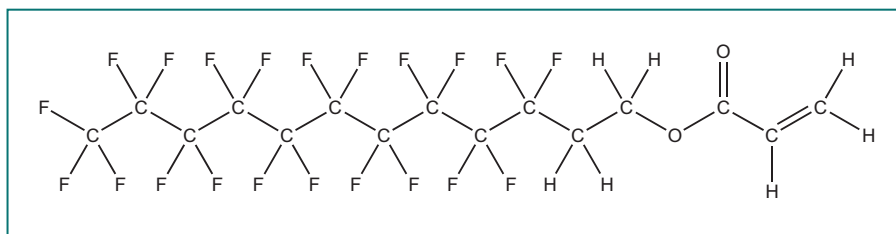
| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|------------------|---|-------------------|
| 8:2FTAcr | 1H,1H,2H,2H-Perfluorodecyl acrylate | 1.2 ml 50 µg/ml |
| 10:2FTAcr | 1H,1H,2H,2H-Perfluorododecyl acrylate (97%) | 1.2 ml 48.5 µg/ml |

NATIVE TELOMER ACETATES (X:2FTOAc)

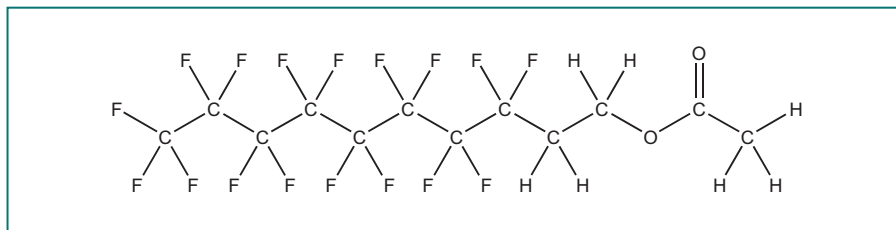
| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|------------------|--------------------------------------|-----------------|
| 8:2FTOAc | 1H,1H,2H,2H-Perfluorodecyl acetate | 1.2 ml 50 µg/ml |
| 10:2FTOAc | 1H,1H,2H,2H-Perfluorododecyl acetate | 1.2 ml 50 µg/ml |



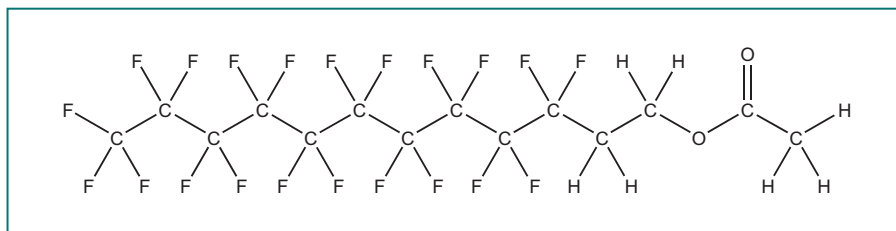
8:2FTAcr



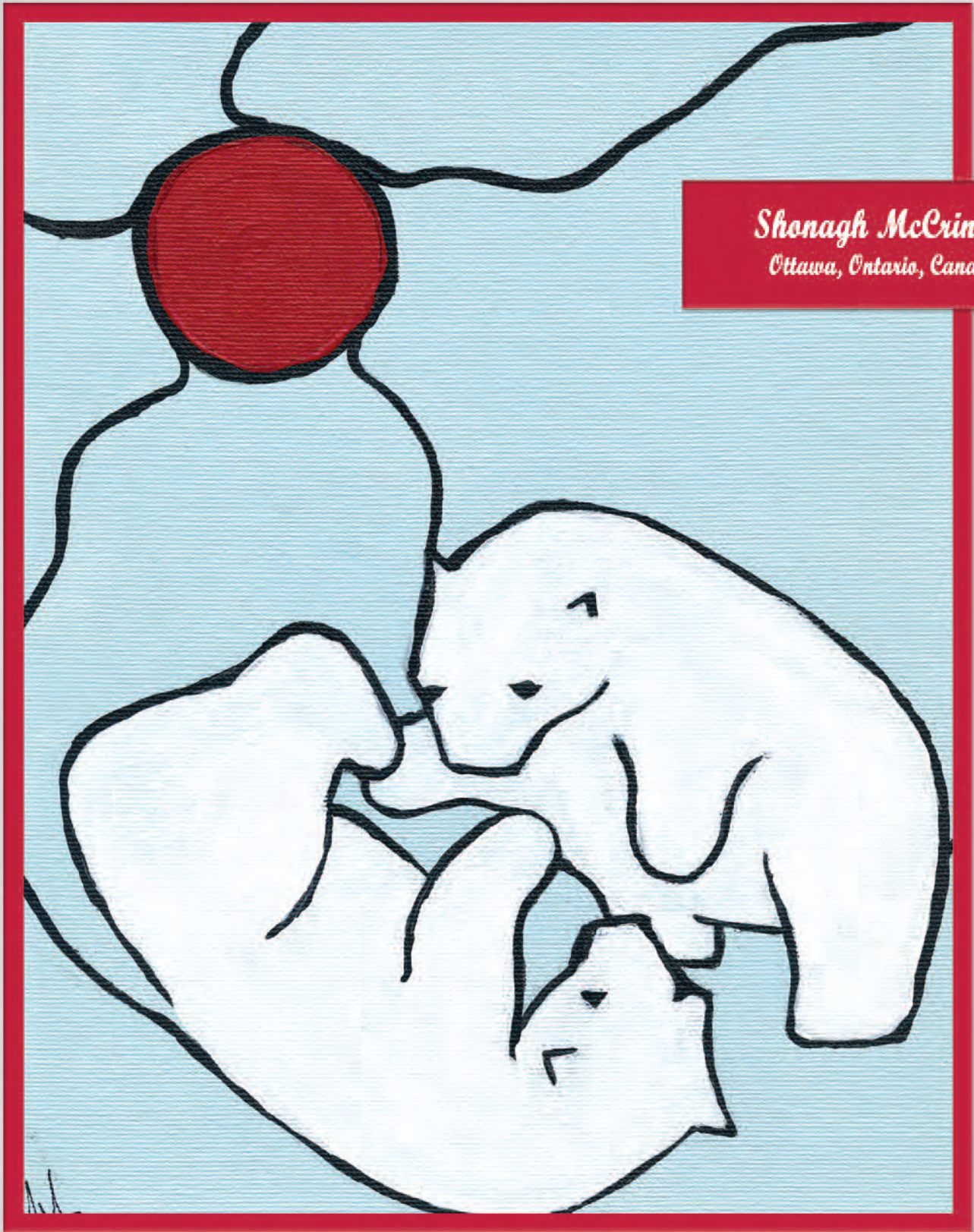
10:2FTAcr



8:2FTOAc



10:2FTOAc



Shonagh McCrindle
Ottawa, Ontario, Canada

CERTIFIED REFERENCE MATERIALS

Wellington currently offers three certified reference materials (CRMs) for use in testing an analytical lab's ability to generate accurate and reproducible data using real, as opposed to fortified, samples.

The following CRMs are currently offered:

WMS-01: Lake Sediment for Organic Contaminant Analysis.

WMF-01: Freeze-Dried Fish Tissue for Organic Contaminant Analysis.

CARP-2: Fish Tissue for Organic Contaminant Analysis.

More details on each of these CRMs, including the analytes and their certified values, are given in the following section.



WMS-01: REFERENCE LAKE SEDIMENT for ORGANIC CONTAMINANT ANALYSIS

| Catalogue Number | Product | Qty/Conc |
|--|---------------------------------|---|
| WMS-01 | Reference Lake Sediment, WMS-01 | 25 g |
| Chlorinated Dibenzo-p-dioxins (PCDDs) | | Certified Reference value (pg/g) |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin Total Tetrachlorodibenzo-p-dioxins | | 17.7 ± 5.6 60.1 ± 25 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin Total Pentachlorodibenzo-p-dioxins | | 7.96 ± 2.8 69.5 ± 23 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin Total Hexachlorodibenzo-p-dioxins | | 8.66 ± 2.7 20.8 ± 4.8 17.3 ± 8.0 238 ± 86 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin Total Heptachlorodibenzo-p-dioxins | | 293 ± 63 608 ± 152 |
| Octachlorodibenzo-p-dioxin | | 1899 ± 456 |
| Chlorinated Dibenzofurans (PCDFs) | | |
| 2,3,7,8-Tetrachlorodibenzofuran Total Tetrachlorodibenzofurans | | 52.5 ± 16 374 ± 162 |
| 1,2,3,7,8-Pentachlorodibenzofuran 2,3,4,7,8-Pentachlorodibenzofuran Total Pentachlorodibenzofurans | | 12.6 ± 5.0 18.5 ± 6.1 225 ± 113 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran 1,2,3,6,7,8-Hexachlorodibenzofuran 1,2,3,7,8,9-Hexachlorodibenzofuran 2,3,4,6,7,8-Hexachlorodibenzofuran Total Hexachlorodibenzofurans | | 67.3 ± 24 20.3 ± 8.7 2.68* ± 4.0 16 ± 8.0 262 ± 95 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran 1,2,3,4,7,8,9-Heptachlorodibenzofuran Total Heptachlorodibenzofurans | | 299 ± 73 15.1 ± 4.6 411 ± 100 |
| Octachlorodibenzofuran | | 509 ± 157 |
| Chlorinated Biphenyls (PCBs-IUPAC) | | |
| 3,3',4,4'-Tetrachlorobiphenyl (77) | | 1717 ± 520 |
| 3,4,4',5-Tetrachlorobiphenyl (81) | | 75* ± 79 |
| 2,3,3',4,4'-Pentachlorobiphenyl (105) | | 3998 ± 951 |
| 2,3,4,4',5-Pentachlorobiphenyl (114) | | 207 ± 128 |
| 2,3',4,4',5-Pentachlorobiphenyl (118) | | 8115 ± 1663 |
| 2',3,4,4',5-Pentachlorobiphenyl (123) | | 209 ± 191 |
| 3,3',4,4',5-Pentachlorobiphenyl (126) | | 84.9 ± 35 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (156) | | 715 ± 248 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl (157) | | 186 ± 81 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl (167) | | 330 ± 85 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl (169) | | 7.97 ± 5.3 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl (189) | | 85.2 ± 17.8 |

* Provisional value for information purposes only. Any negative deviation is inadmissible.
The concentrations of these analytes may be certified at a later date as more data becomes available.

WMF-01: REFERENCE FISH TISSUE for ORGANIC CONTAMINANT ANALYSIS

| Catalogue Number | Product | Qty/Conc |
|---|--|----------------|
| WMF-01 | Reference "Freeze-Dried" Fish Tissue, WMF-01 | 1 x 10 g |
| Chlorinated Dibenzo-p-dioxins (PCDDs) | | |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | | 13.1 ± 4.4 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | | 2.72 ± 1.3 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | | 0.22* ± 0.3 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | | 0.88 ± 0.4 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | | 0.27* ± 0.4 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | | 0.59* ± 0.7 |
| Octachlorodibenzo-p-dioxin | | 3.91* ± 6.2 |
| Chlorinated Dibenzofurans (PCDFs) | | |
| 2,3,7,8-Tetrachlorodibenzofuran | | 13.1 ± 4.9 |
| 1,2,3,7,8-Pentachlorodibenzofuran | | 1.53* ± 1.4 |
| 2,3,4,7,8-Pentachlorodibenzofuran | | 7.15 ± 2.2 |
| 1,2,3,4,7,8-Hexachlorodibenzofuran | | 0.86* ± 1.0 |
| 1,2,3,6,7,8-Hexachlorodibenzofuran | | 0.51* ± 0.7 |
| 1,2,3,7,8,9-Hexachlorodibenzofuran | | 0.25* ± 0.4 |
| 2,3,4,6,7,8-Hexachlorodibenzofuran | | 0.68* ± 1.2 |
| 1,2,3,4,6,7,8-Heptachlorodibenzofuran | | 1.01* ± 1.9 |
| 1,2,3,4,7,8,9-Heptachlorodibenzofuran | | 0.30* ± 0.5 |
| Octachlorodibenzofuran | | 1.38* ± 2.1 |
| Chlorinated Biphenyls (PCBs-IUPAC) | | |
| 3,3',4,4'-Tetrachlorobiphenyl (77) | | 2233 ± 720 |
| 3,4,4',5-Tetrachlorobiphenyl (81) | | 201 ± 58 |
| 2,3,3',4,4'-Pentachlorobiphenyl (105) | | 49050 ± 14200 |
| 2,3,4,4',5-Pentachlorobiphenyl (114) | | 3523 ± 1670 |
| 2,3',4,4',5-Pentachlorobiphenyl (118) | | 130100 ± 32500 |
| 2',3,4,4',5-Pentachlorobiphenyl (123) | | 4233* ± 2620 |
| 3,3',4,4',5-Pentachlorobiphenyl (126) | | 739 ± 260 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (156) | | 14890 ± 5020 |
| 2,3,3',4,4',5'-Hexachlorobiphenyl (157) | | 3488 ± 870 |
| 2,3',4,4',5,5'-Hexachlorobiphenyl (167) | | 9750 ± 3090 |
| 3,3',4,4',5,5'-Hexachlorobiphenyl (169) | | 76 ± 30 |
| 2,3,3',4,4',5,5'-Heptachlorobiphenyl (189) | | 2016 ± 611 |
| Brominated Diphenyl Ethers (PBDEs-IUPAC) | | |
| 2,4,4'-Tribromodiphenyl ether (28) | | 3124 ± 290 |
| 2,2',4,4'-Tetrabromodiphenyl ether (47) | | 123200 ± 24800 |
| 2,2',4,4',5-Pentabromodiphenyl ether (99) | | 37500 ± 4220 |
| 2,2',4,4',6-Pentabromodiphenyl ether (100) | | 35870 ± 14500 |
| 2,2',4,4',5,5'-Hexabromodiphenyl ether (153) | | 17040 ± 8000 |
| 2,2',4,4',5,6'-Hexabromodiphenyl ether (154) | | 19790 ± 2880 |
| 2,2',3,4,4',5',6-Heptabromodiphenyl ether (183) | | 532* ± 400 |

* Provisional value for information purposes only. Any negative deviation is inadmissible.
The concentrations of these analytes may be certified at a later date as more data becomes available.

CARP-2: REFERENCE FISH TISSUE for ORGANIC CONTAMINANT ANALYSIS

| Catalogue Number | Product | Qty/Conc |
|---|-------------------------------|--|
| CARP-2 | Reference Fish Tissue, CARP-2 | 6 x 9 g |
| Polychlorinated Biphenyls (PCBs) Congener (IUPAC) | | Certified Concentration µg/kg (wet weight basis) |
| 18 | | 27.3 ± 4.0 |
| 28 | | 34.0 ± 7.2 |
| 44 | | 86.6 ± 25.9 |
| 52 | | 138 ± 43 |
| 118 | | 148 ± 33 |
| 128 | | 20.4 ± 4.4 |
| 153 | | 105 ± 22 |
| 180 | | 53.3 ± 13.0 |
| 194 | | 10.9 ± 3.1 |
| 206 | | 4.4 ± 1.1 |
| Polychlorinated Biphenyls (PCBs) Congener (IUPAC) | | Reference Concentration* µg/kg (wet weight basis) |
| 8 | | 4.8 ± 1.8 |
| 66/95 | | 174 ± 52 |
| 101/90 | | 145 ± 48 |
| 105 | | 53.2 ± 15.6 |
| 138/163/164 | | 103 ± 30 |
| 170/190 | | 20.6 ± 2.9 |
| 187/182 | | 37.1 ± 6.3 |
| 209 | | 4.6 ± 2.0 |
| Polychlorinated dibenzo-p-dioxins (PCDDs) | | ng/kg (wet weight basis)* |
| 2,3,7,8-Tetrachlorodibenzo-p-dioxin | | 7.4 ± 0.7 |
| 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | | 5.3 ± 1.3 |
| 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | | 1.6 ± 0.3 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | | 5.8 ± 0.8 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | | 0.78 ± 0.12 |
| 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | | 6.4 ± 0.9 |
| Octachlorodibenzo-p-dioxin | | 9.4 ± 1.7 |
| Polychlorinated dibenzofurans (PCDFs) | | ng/kg (wet weight basis)* |
| 2,3,7,8-Tetrachlorodibenzofuran | | 18.2 ± 1.6 |
| 1,2,3,7,8-Pentachlorodibenzofuran | | 5.6 ± 0.3 |
| Pesticides | | µg/kg (wet weight basis)* |
| gamma-chlordane | | 4.5 ± 0.7 |
| 2,4'-DDE | | 2.9 ± 0.5 |
| trans-nonachlor | | 11.0 ± 0.9 |
| dieldrin | | 8.3 ± 0.8 |
| 4,4'-DDE | | 158 ± 14 |
| 2,4'-DDD | | 21.8 ± 0.7 |
| 4,4'-DDD | | 90.9 ± 8.5 |

* Not Certified

CARP-2 was prepared and certified by the National Research Council of Canada (NRCC), Institute for Environmental Research and Technology.

ADDITIONAL PRODUCTS

PAH Calibration Sets

and Native and Mass-Labelled Support Solutions

Chlorinated Biphenyls (HO-PCBs):

Native and Mass-Labelled Individuals;
Mass-Labelled Solution/Mixture

Methoxy Chlorobiphenyls (MeO-PCBs):

Native and Mass-Labelled Individuals and Solution/Mixtures

Chlorinated Biphenylenes (PCBPs):

Native and Mass-Labelled

Triclocarban:

Native and Mass-Labelled

Triclosan and Methyl Triclosan:

Native and Mass-Labelled and Chlorinated Derivatives

Tris(4-chlorophenyl) Methane and Methanol:

Native and Mass-Labelled

Chlorinated Naphthalenes (PCNs):

Native Individuals and Solution/Mixtures

Chlorinated Diphenyl Ethers (PCDEs):

Native and Mass-Labelled

Native and Mass-Labelled Chlorobenzene and Chlorophenol Solution/Mixtures

Mass-Labelled Chlorobenzenes and Chlorophenols

Melamine and Cyanuric Acid:

Native and Mass-Labelled

Native and Mass-Labelled Bisphenol A and Native Bisphenol Analogues

Native and Mass-Labelled Tetrachlorodibenzothiophenes

Native and Mass-Labelled Halogenated Carbazoles



PAH-CVS-A

| Catalogue Number | Product (isooctane/toluene solution) | Qty/Conc |
|------------------|--|-----------------------|
| PAH-CVS-A | PAH-CVS-A Calibration Solutions CS1-CS5 | 1 kit (5 ampoules) |
| PAH-A-CS1 | CS1 | 1.0 ml |
| PAH-A-CS2 | CS2 | 1.0 ml |
| PAH-A-CS3 | CS3 | 1.0 ml |
| PAH-A-CS4 | CS4 | 1.0 ml |
| PAH-A-CS5 | CS5 | 1.0 ml |

| | PAH-A- CS1 (ng/ml) | PAH-A- CS2 (ng/ml) | PAH-A- CS3 (ng/ml) | PAH-A- CS4 (ng/ml) | PAH-A- CS5 (ng/ml) |
|--|--------------------------|--------------------------|--------------------------|--------------------------|--------------------------|
| Native PAHs | | | | | |
| Naphthalene | 2.0 | 10 | 40 | 200 | 800 |
| Acenaphthylene | 2.0 | 10 | 40 | 200 | 800 |
| Acenaphthene | 2.0 | 10 | 40 | 200 | 800 |
| Fluorene | 2.0 | 10 | 40 | 200 | 800 |
| Phenanthrene | 2.0 | 10 | 40 | 200 | 800 |
| Anthracene | 2.0 | 10 | 40 | 200 | 800 |
| Fluoranthene | 2.0 | 10 | 40 | 200 | 800 |
| Pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Benz[a]anthracene | 2.0 | 10 | 40 | 200 | 800 |
| Cyclopenta[c,d]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Chrysene | 2.0 | 10 | 40 | 200 | 800 |
| 5-Methylchrysene | 2.0 | 10 | 40 | 200 | 800 |
| Benzo[b]fluoranthene | 2.0 | 10 | 40 | 200 | 800 |
| Benzo[k]fluoranthene | 2.0 | 10 | 40 | 200 | 800 |
| Benzo[j]fluoranthene | 2.0 | 10 | 40 | 200 | 800 |
| Benzo[a]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Indeno[1,2,3-c,d]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Dibenz[a,h]anthracene | 2.0 | 10 | 40 | 200 | 800 |
| Benzo[g,h,i]perylene | 2.0 | 10 | 40 | 200 | 800 |
| Dibenzo[a,l]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Dibenzo[a,e]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Dibenzo[a,i]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Dibenzo[a,h]pyrene | 2.0 | 10 | 40 | 200 | 800 |
| Deuterated PAHs | | | | | |
| Naphthalene-d ₈ | 100 | 100 | 100 | 100 | 100 |
| Acenaphthene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Fluorene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Phenanthrene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Anthracene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Fluoranthene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Pyrene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Benz[a]anthracene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Chrysene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[b]fluoranthene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[k]fluoranthene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[a]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Dibenz[a,h]anthracene-d ₁₄ | 100 | 100 | 100 | 100 | 100 |
| Benzo[g,h,i]perylene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Dibenzo[a,i]pyrene-d ₁₄ | 100 | 100 | 100 | 100 | 100 |
| Deuterated PAH Internal Standards | | | | | |
| Acenaphthylene-d ₈ | 100 | 100 | 100 | 100 | 100 |
| p-Terphenyl-d ₁₄ | 100 | 100 | 100 | 100 | 100 |
| Benzo[e]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |

| Catalogue Number | Product (isooctane/toluene solution) | Qty/Conc |
|------------------|--|----------|
| PAH-LCS-A | PAH Labelled Compound Solution | 1.2 ml |
| PAH-ISS-A | PAH Internal Standard Spiking Solution | 1.2 ml |
| PAH-STK-A | PAH Native Stock Solution | 1.2 ml |

| | PAH-LCS-A (ng/ml) | PAH-ISS-A (ng/ml) | PAH-STK-A (ng/ml) |
|--|----------------------|----------------------|----------------------|
| Native PAHs | | | |
| Naphthalene | — | — | 4000 |
| Acenaphthylene | — | — | 4000 |
| Acenaphthene | — | — | 4000 |
| Fluorene | — | — | 4000 |
| Phenanthrene | — | — | 4000 |
| Anthracene | — | — | 4000 |
| Fluoranthene | — | — | 4000 |
| Pyrene | — | — | 4000 |
| Benz[a]anthracene | — | — | 4000 |
| Cyclopenta[c,d]pyrene | — | — | 4000 |
| Chrysene | — | — | 4000 |
| 5-Methylchrysene | — | — | 4000 |
| Benzo[b]fluoranthene | — | — | 4000 |
| Benzo[k]fluoranthene | — | — | 4000 |
| Benzo[j]fluoranthene | — | — | 4000 |
| Benzo[a]pyrene | — | — | 4000 |
| Indeno[1,2,3-c,d]pyrene | — | — | 4000 |
| Dibenz[a,h]anthracene | — | — | 4000 |
| Benzo[g,h,i]perylene | — | — | 4000 |
| Dibenzo[a,l]pyrene | — | — | 4000 |
| Dibenzo[a,e]pyrene | — | — | 4000 |
| Dibenzo[a,i]pyrene | — | — | 4000 |
| Dibenzo[a,h]pyrene | — | — | 4000 |
| Deuterated PAHs | | | |
| Naphthalene-d ₈ | 2000 | — | — |
| Acenaphthene-d ₁₀ | 2000 | — | — |
| Fluorene-d ₁₀ | 2000 | — | — |
| Phenanthrene-d ₁₀ | 2000 | — | — |
| Anthracene-d ₁₀ | 2000 | — | — |
| Fluoranthene-d ₁₀ | 2000 | — | — |
| Pyrene-d ₁₀ | 2000 | — | — |
| Benz[a]anthracene-d ₁₂ | 2000 | — | — |
| Chrysene-d ₁₂ | 2000 | — | — |
| Benzo[b]fluoranthene-d ₁₂ | 2000 | — | — |
| Benzo[k]fluoranthene-d ₁₂ | 2000 | — | — |
| Benzo[a]pyrene-d ₁₂ | 2000 | — | — |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | 2000 | — | — |
| Dibenz[a,h]anthracene-d ₁₄ | 2000 | — | — |
| Benzo[g,h,i]perylene-d ₁₂ | 2000 | — | — |
| Dibenzo[a,i]pyrene-d ₁₄ | 2000 | — | — |
| Deuterated PAH Internal Standards | | | |
| Acenaphthylene-d ₈ | — | 2000 | — |
| p-Terphenyl-d ₁₄ | — | 2000 | — |
| Benzo[e]pyrene-d ₁₂ | — | 2000 | — |

PAH-CVS-B

| Catalogue Number | Product (isooctane/toluene solution) | Qty/Conc |
|------------------|--------------------------------------|--------------|
| PAH-CVS-B | PAH-CVS-B | 1 kit |
| | Calibration Solutions CS1-CS5 | (5 ampoules) |
| PAH-B-CS1 | CS1 | 1.0 ml |
| PAH-B-CS2 | CS2 | 1.0 ml |
| PAH-B-CS3 | CS3 | 1.0 ml |
| PAH-B-CS4 | CS4 | 1.0 ml |
| PAH-B-CS5 | CS5 | 1.0 ml |

| | PAH-B-CS1 (ng/ml) | PAH-B-CS2 (ng/ml) | PAH-B-CS3 (ng/ml) | PAH-B-CS4 (ng/ml) | PAH-B-CS5 (ng/ml) |
|---|----------------------|----------------------|----------------------|----------------------|----------------------|
| Native PAHs | | | | | |
| Naphthalene | 2.0 | 10 | 50 | 250 | 1000 |
| 2-Methylnaphthalene | 2.0 | 10 | 50 | 250 | 1000 |
| Acenaphthylene | 2.0 | 10 | 50 | 250 | 1000 |
| Acenaphthene | 2.0 | 10 | 50 | 250 | 1000 |
| Fluorene | 2.0 | 10 | 50 | 250 | 1000 |
| Phenanthrene | 2.0 | 10 | 50 | 250 | 1000 |
| Anthracene | 2.0 | 10 | 50 | 250 | 1000 |
| Fluoranthene | 2.0 | 10 | 50 | 250 | 1000 |
| Pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[c]fluorene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[a]anthracene | 2.0 | 10 | 50 | 250 | 1000 |
| Cyclopenta[c,d]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Chrysene | 2.0 | 10 | 50 | 250 | 1000 |
| 5-Methylchrysene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[b]fluoranthene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[k]fluoranthene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[j]fluoranthene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[e]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[a]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Perylene | 2.0 | 10 | 50 | 250 | 1000 |
| Indeno[1,2,3-c,d]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Dibenz[a,h]anthracene | 2.0 | 10 | 50 | 250 | 1000 |
| Benzo[g,h,i]perylene | 2.0 | 10 | 50 | 250 | 1000 |
| Dibenzo[a,l]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Dibenzo[a,e]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Dibenzo[a,i]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Dibenzo[a,h]pyrene | 2.0 | 10 | 50 | 250 | 1000 |
| Deuterated PAHs (PAH-LCS-B) | | | | | |
| Naphthalene-d ₈ | 100 | 100 | 100 | 100 | 100 |
| 2-Methylnaphthalene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Acenaphthylene-d ₈ | 100 | 100 | 100 | 100 | 100 |
| Phenanthrene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Anthracene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Fluoranthene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Benzo[a]anthracene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Chrysene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[b]fluoranthene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[k]fluoranthene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Benzo[a]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Perylene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Dibenz[a,h]anthracene-d ₁₄ | 100 | 100 | 100 | 100 | 100 |
| Benzo[g,h,i]perylene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Dibenzo[a,i]pyrene-d ₁₄ | 100 | 100 | 100 | 100 | 100 |
| Internal Standards (PAH-ISS-B) | | | | | |
| Acenaphthene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Pyrene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| Benzo[e]pyrene-d ₁₂ | 100 | 100 | 100 | 100 | 100 |
| Sampling Standards (PAH-SS-B) | | | | | |
| Fluorene-d ₁₀ | 100 | 100 | 100 | 100 | 100 |
| p-Terphenyl-d ₁₄ | 100 | 100 | 100 | 100 | 100 |

| Catalogue Number | Product (isooctane/toluene solution) | | | Qty/Conc |
|---|--|------------------------------|-----------------------------|------------------------------|
| PAH-LCS-B | PAH Labelled Compound Solution | | | 1.2 ml |
| PAH-ISS-B | PAH Internal Standard Spiking Solution | | | 1.2 ml |
| PAH-SS-B | PAH Sampling Standard Solution | | | 1.2 ml |
| PAH-STK-B | PAH Native Stock Solution | | | 1.2 ml |
| | PAH-LCS-B (ng/ml) | PAH-ISS-B (ng/ml) | PAH-SS-B (ng/ml) | PAH-STK-B (ng/ml) |
| Native PAHs | | | | |
| Naphthalene | — | — | — | 2500 |
| 2-Methylnaphthalene | — | — | — | 2500 |
| Acenaphthylene | — | — | — | 2500 |
| Acenaphthene | — | — | — | 2500 |
| Fluorene | — | — | — | 2500 |
| Phenanthrene | — | — | — | 2500 |
| Anthracene | — | — | — | 2500 |
| Fluoranthene | — | — | — | 2500 |
| Pyrene | — | — | — | 2500 |
| Benzo[c]fluorene | — | — | — | 2500 |
| Benz[a]anthracene | — | — | — | 2500 |
| Cyclopenta[c,d]pyrene | — | — | — | 2500 |
| Chrysene | — | — | — | 2500 |
| 5-Methylchrysene | — | — | — | 2500 |
| Benzo[b]fluoranthene | — | — | — | 2500 |
| Benzo[k]fluoranthene | — | — | — | 2500 |
| Benzo[j]fluoranthene | — | — | — | 2500 |
| Benzo[e]pyrene | — | — | — | 2500 |
| Benzo[a]pyrene | — | — | — | 2500 |
| Perylene | — | — | — | 2500 |
| Indeno[1,2,3-c,d]pyrene | — | — | — | 2500 |
| Dibenz[a,h]anthracene | — | — | — | 2500 |
| Benzo[g,h,i]perylene | — | — | — | 2500 |
| Dibenzo[a,i]pyrene | — | — | — | 2500 |
| Dibenzo[a,e]pyrene | — | — | — | 2500 |
| Dibenzo[a,i]pyrene | — | — | — | 2500 |
| Dibenzo[a,h]pyrene | — | — | — | 2500 |
| Deuterated PAHs | | | | |
| Naphthalene-d ₈ | 5000 | — | — | — |
| 2-Methylnaphthalene-d ₁₀ | 5000 | — | — | — |
| Acenaphthylene-d ₈ | 5000 | — | — | — |
| Phenanthrene-d ₁₀ | 5000 | — | — | — |
| Anthracene-d ₁₀ | 5000 | — | — | — |
| Fluoranthene-d ₁₀ | 5000 | — | — | — |
| Benz[a]anthracene-d ₁₂ | 5000 | — | — | — |
| Chrysene-d ₁₂ | 5000 | — | — | — |
| Benzo[b]fluoranthene-d ₁₂ | 5000 | — | — | — |
| Benzo[k]fluoranthene-d ₁₂ | 5000 | — | — | — |
| Benzo[a]pyrene-d ₁₂ | 5000 | — | — | — |
| Perylene-d ₁₂ | 5000 | — | — | — |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | 5000 | — | — | — |
| Dibenz[a,h]anthracene-d ₁₄ | 5000 | — | — | — |
| Benzo[g,h,i]perylene-d ₁₂ | 5000 | — | — | — |
| Dibenzo[a,i]pyrene-d ₁₄ | 5000 | — | — | — |
| Internal Standards | | | | |
| Acenaphthene-d ₁₀ | — | 5000 | — | — |
| Pyrene-d ₁₀ | — | 5000 | — | — |
| Benzo[e]pyrene-d ₁₂ | — | 5000 | — | — |
| Sampling Standards | | | | |
| Fluorene-d ₁₀ | — | — | 5000 | — |
| p-Terphenyl-d ₁₄ | — | — | 5000 | — |

METHOD 429: HRGC/LRMS CALIBRATION SOLUTIONS FOR PAHs

| Catalogue Number | Product (isooctane/toluene solution) | Qty/Conc |
|------------------|---|-----------------------|
| L429-CVS | L429-CVS Calibration Solutions CS1-CS5 | 1 kit (5 ampoules) |
| L429-CS1 | CS1 | 1.0 ml |
| L429-CS2 | CS2 | 1.0 ml |
| L429-CS3 | CS3 | 1.0 ml |
| L429-CS4 | CS4 | 1.0 ml |
| L429-CS5 | CS5 | 1.0 ml |

| | L429-CS1 (ng/μl) | L429-CS2 (ng/μl) | L429-CS3 (ng/μl) | L429-CS4 (ng/μl) | L429-CS5 (ng/μl) |
|---|---------------------|---------------------|---------------------|---------------------|---------------------|
| Native PAHs | | | | | |
| Naphthalene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| 2-Methylnaphthalene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Acenaphthylene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Acenaphthene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Fluorene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Phenanthrene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Anthracene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Fluoranthene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Pyrene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benz[a]anthracene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Chrysene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benzo[b]fluoranthene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benzo[k]fluoranthene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benzo[e]pyrene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benzo[a]pyrene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Perylene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Indeno[1,2,3-c,d]pyrene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Dibenz[a,h]anthracene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Benzo[g,h,i]perylene | 0.25 | 0.5 | 1.0 | 2.5 | 5.0 |
| Surrogate Standards | | | | | |
| Fluorene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| p-Terphenyl-d ₁₄ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Internal Standards | | | | | |
| Naphthalene-d ₈ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| 2-Methylnaphthalene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Acenaphthylene-d ₈ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Phenanthrene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Fluoranthene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[a]anthracene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Chrysene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[b]fluoranthene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[k]fluoranthene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[a]pyrene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Perylene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Dibenz[a,h]anthracene-d ₁₄ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[g,h,i]perylene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Alternate Standard | | | | | |
| Anthracene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Recovery Standards | | | | | |
| Acenaphthene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Pyrene-d ₁₀ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| Benzo[e]pyrene-d ₁₂ | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |

METHOD 429: HRGC/LRMS CALIBRATION SOLUTIONS FOR PAHs

| Catalogue Number | Product (isooctane/toluene solution) | | | | | Qty/Conc |
|------------------|--|--|--|--|--|----------|
| L429-SS | Method 429 Surrogate Standard Stock Solution | | | | | 1.2 ml |
| L429-IS | Method 429 Internal Standard Stock Solution | | | | | 1.2 ml |
| L429-AS | Method 429 Alternate Standard Stock Solution | | | | | 1.2 ml |
| L429-RS | Method 429 Recovery Standard Stock Solution | | | | | 1.2 ml |
| L429-PAR | Method 429 Native PAH Stock Solution | | | | | 1.2 ml |

| | L429-SS (µg/ml) | L429-IS (µg/ml) | L429-AS (µg/ml) | L429-RS (µg/ml) | L429-PAR (µg/ml) |
|---|--------------------|--------------------|--------------------|--------------------|---------------------|
| Native PAHs | | | | | |
| Naphthalene | — | — | — | — | 2.0 |
| 2-Methylnaphthalene | — | — | — | — | 2.0 |
| Acenaphthylene | — | — | — | — | 2.0 |
| Acenaphthene | — | — | — | — | 2.0 |
| Fluorene | — | — | — | — | 2.0 |
| Phenanthrene | — | — | — | — | 2.0 |
| Anthracene | — | — | — | — | 2.0 |
| Fluoranthene | — | — | — | — | 2.0 |
| Pyrene | — | — | — | — | 2.0 |
| Benz[a]anthracene | — | — | — | — | 2.0 |
| Chrysene | — | — | — | — | 2.0 |
| Benzo[b]fluoranthene | — | — | — | — | 2.0 |
| Benzo[k]fluoranthene | — | — | — | — | 2.0 |
| Benzo[e]pyrene | — | — | — | — | 2.0 |
| Benzo[a]pyrene | — | — | — | — | 2.0 |
| Perylene | — | — | — | — | 2.0 |
| Indeno[1,2,3-c,d]pyrene | — | — | — | — | 2.0 |
| Dibenz[a,h]anthracene | — | — | — | — | 2.0 |
| Benzo[g,h,i]perylene | — | — | — | — | 2.0 |
| Surrogate Standards | | | | | |
| Fluorene-d ₁₀ | 100 | — | — | — | — |
| p-Terphenyl-d ₁₄ | 100 | — | — | — | — |
| Internal Standards | | | | | |
| Naphthalene-d ₈ | — | 100 | — | — | — |
| 2-Methylnaphthalene-d ₁₀ | — | 100 | — | — | — |
| Acenaphthylene-d ₈ | — | 100 | — | — | — |
| Phenanthrene-d ₁₀ | — | 100 | — | — | — |
| Fluoranthene-d ₁₀ | — | 100 | — | — | — |
| Benz[a]anthracene-d ₁₂ | — | 100 | — | — | — |
| Chrysene-d ₁₂ | — | 100 | — | — | — |
| Benzo[b]fluoranthene-d ₁₂ | — | 100 | — | — | — |
| Benzo[k]fluoranthene-d ₁₂ | — | 100 | — | — | — |
| Benzo[a]pyrene-d ₁₂ | — | 100 | — | — | — |
| Perylene-d ₁₂ | — | 100 | — | — | — |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | — | 100 | — | — | — |
| Dibenz[a,h]anthracene-d ₁₄ | — | 100 | — | — | — |
| Benzo[g,h,i]perylene-d ₁₂ | — | 100 | — | — | — |
| Alternate Standard | | | | | |
| Anthracene-d ₁₀ | — | — | 100 | — | — |
| Recovery Standards | | | | | |
| Acenaphthene-d ₁₀ | — | — | — | 100 | — |
| Pyrene-d ₁₀ | — | — | — | 100 | — |
| Benzo[e]pyrene-d ₁₂ | — | — | — | 100 | — |

EPA PAH SOLUTION/MIXTURES

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|--------------------|-----------------------------|----------|
| EPA-PAH-STK | Native PAH Solution/Mixture | 1.2 ml |
| | Naphthalene | 5 µg/ml |
| | Acenaphthylene | 5 µg/ml |
| | Acenaphthene | 5 µg/ml |
| | Fluorene | 5 µg/ml |
| | Phenanthrene | 5 µg/ml |
| | Anthracene | 5 µg/ml |
| | Fluoranthene | 5 µg/ml |
| | Pyrene | 5 µg/ml |
| | Benzo[a]anthracene | 5 µg/ml |
| | Chrysene | 5 µg/ml |
| | Benzo[b]fluoranthene | 5 µg/ml |
| | Benzo[k]fluoranthene | 5 µg/ml |
| | Benzo[a]pyrene | 5 µg/ml |
| | Indeno[1,2,3-c,d]pyrene | 5 µg/ml |
| | Benzo[g,h,i]perylene | 5 µg/ml |
| | Dibenzo[a,h]anthracene | 5 µg/ml |

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|--------------------|---|----------|
| EPA-PAH-LCS | Deuterated PAH Solution/Mixture | 1.2 ml |
| | Naphthalene-d ₈ | 5 µg/ml |
| | Acenaphthylene-d ₈ | 5 µg/ml |
| | Acenaphthene-d ₁₀ | 5 µg/ml |
| | Fluorene-d ₁₀ | 5 µg/ml |
| | Phenanthrene-d ₁₀ | 5 µg/ml |
| | Anthracene-d ₁₀ | 5 µg/ml |
| | Fluoranthene-d ₁₀ | 5 µg/ml |
| | Pyrene-d ₁₀ | 5 µg/ml |
| | Benzo[a]anthracene-d ₁₂ | 5 µg/ml |
| | Chrysene-d ₁₂ | 5 µg/ml |
| | Benzo[b]fluoranthene-d ₁₂ | 5 µg/ml |
| | Benzo[k]fluoranthene-d ₁₂ | 5 µg/ml |
| | Benzo[a]pyrene-d ₁₂ | 5 µg/ml |
| | Indeno[1,2,3-c,d]pyrene-d ₁₂ | 5 µg/ml |
| | Benzo[g,h,i]perylene-d ₁₂ | 5 µg/ml |
| | Dibenzo[a,h]anthracene-d ₁₄ | 5 µg/ml |

* The solutions above can be used with the **PAH-CVS-B** calibration solutions.

EPA & EU PAH SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|-------------------------------------|---------------------------------|----------|
| EPA-EU-PAH-ISS | Deuterated PAH Solution/Mixture | 1.2 ml |
| 2-Methylnaphthalene-d ₁₀ | | 5 µg/ml |
| p-Terphenyl-d ₁₄ | | 5 µg/ml |
| Benzo[e]pyrene-d ₁₂ | | 5 µg/ml |

EU PAH SOLUTION/MIXTURES

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|-------------------------|-----------------------------|----------|
| EU-PAH-STK | Native PAH Solution/Mixture | 1.2 ml |
| Benzo[c]fluorene | | 5 µg/ml |
| Cyclopenta[c,d]pyrene | | 5 µg/ml |
| Benzo[a]anthracene | | 5 µg/ml |
| Chrysene | | 5 µg/ml |
| 5-Methylchrysene | | 5 µg/ml |
| Benzo[b]fluoranthene | | 5 µg/ml |
| Benzo[k]fluoranthene | | 5 µg/ml |
| Benzo[j]fluoranthene | | 5 µg/ml |
| Benzo[a]pyrene | | 5 µg/ml |
| Indeno[1,2,3-c,d]pyrene | | 5 µg/ml |
| Benzo[g,h,i]perylene | | 5 µg/ml |
| Dibenzo[a,h]anthracene | | 5 µg/ml |
| Dibenzo[a,l]pyrene | | 5 µg/ml |
| Dibenzo[a,e]pyrene | | 5 µg/ml |
| Dibenzo[a,i]pyrene | | 5 µg/ml |
| Dibenzo[a,h]pyrene | | 5 µg/ml |

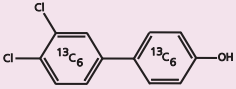
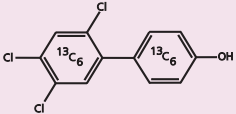
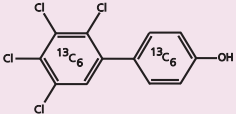
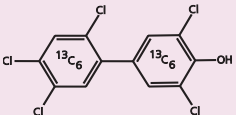
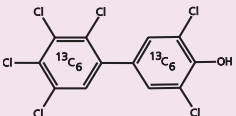
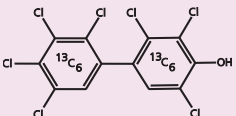
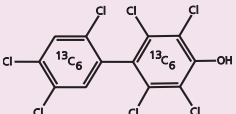
| Catalogue Number | Product (toluene solution) | Qty/Conc |
|---|---------------------------------|----------|
| EU-PAH-LCS | Deuterated PAH Solution/Mixture | 1.2 ml |
| Benzo[a]anthracene-d ₁₂ | | 5 µg/ml |
| Chrysene-d ₁₂ | | 5 µg/ml |
| Benzo[b]fluoranthene-d ₁₂ | | 5 µg/ml |
| Benzo[k]fluoranthene-d ₁₂ | | 5 µg/ml |
| Benzo[a]pyrene-d ₁₂ | | 5 µg/ml |
| Indeno[1,2,3-c,d]pyrene-d ₁₂ | | 5 µg/ml |
| Benzo[g,h,i]perylene-d ₁₂ | | 5 µg/ml |
| Dibenzo[a,h]anthracene-d ₁₄ | | 5 µg/ml |
| Dibenzo[a,i]pyrene-d ₁₄ | | 5 µg/ml |

* The solutions above can be used with the **PAH-CVS-B** calibration solutions.

NATIVE CHLORINATED BIPHENYLS (HO-PCBs)

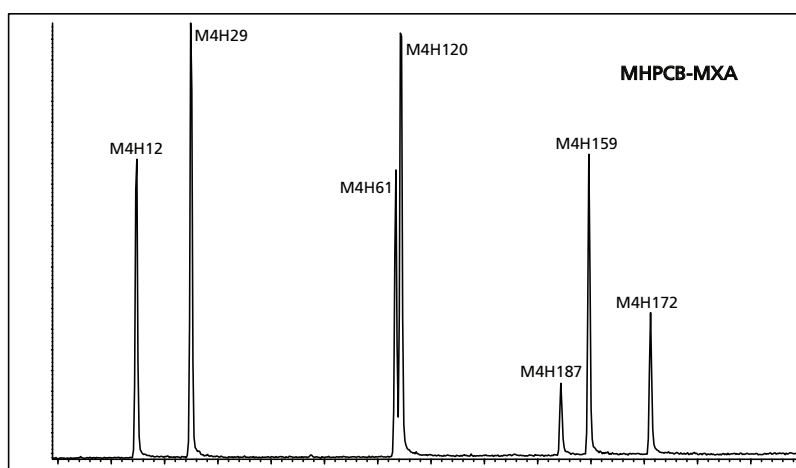
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-------------------|
| 4H107 | 2,3,3',4',5-Pentachloro-4-biphenylol | 1.2 ml 50 µg/ml |
| 4H108 | 2',3,3',4',5-Pentachloro-4-biphenylol (96.5%) | 1.2 ml 48.3 µg/ml |
| 3H118 | 2,3',4,4',5-Pentachloro-3-biphenylol | 1.2 ml 50 µg/ml |
| 4H130 | 2,2',3,3',4',5-Hexachloro-4-biphenylol | 1.2 ml 50 µg/ml |
| 3H138 | 2,2',3',4,4',5-Hexachloro-3-biphenylol | 1.2 ml 50 µg/ml |
| 4H146 | 2,2',3,4',5,5'-Hexachloro-4-biphenylol | 1.2 ml 50 µg/ml |
| 3H153 | 2,2',4,4',5,5'-Hexachloro-3-biphenylol | 1.2 ml 50 µg/ml |
| 4H172 | 2,2',3,3',4',5,5'-Heptachloro-4-biphenylol | 1.2 ml 50 µg/ml |
| 3H180 | 2,2',3',4,4',5,5'-Heptachloro-3-biphenylol | 1.2 ml 50 µg/ml |
| 4H187 | 2,2',3,4',5,5',6-Heptachloro-4-biphenylol | 1.2 ml 50 µg/ml |

MASS-LABELLED CHLORINATED BIPHENYLS

| Catalogue Number | Product |
|------------------|--|
| M4H12 |  <p>3',4'-Dichloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H29 |  <p>2',4',5'-Trichloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H61 |  <p>2',3',4',5'-Tetrachloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H120 |  <p>2',3,4',5,5'-Pentachloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H159 |  <p>2',3,3',4',5,5'-Hexachloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H172 |  <p>2,2',3,3',4',5,5'-Heptachloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |
| M4H187 |  <p>2,2',3,4',5,5',6-Heptachloro-4-[¹³C₁₂]biphenylol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |

MASS-LABELLED CHLORINATED BIPHENYLS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|---|---|----------|
| MHPCB-MXA | Mass-Labelled Chlorinated Biphenyl Solution/Mixture | 1.2 ml |
| 3',4'-Dichloro-4-[¹³ C ₁₂]biphenylol | M4H12 | 5 µg/ml |
| 2',4',5'-Trichloro-4-[¹³ C ₁₂]biphenylol | M4H29 | 5 µg/ml |
| 2',3',4',5'-Tetrachloro-4-[¹³ C ₁₂]biphenylol | M4H61 | 5 µg/ml |
| 2',3,4',5,5'-Pentachloro-4-[¹³ C ₁₂]biphenylol | M4H120 | 5 µg/ml |
| 2',3,3',4',5,5'-Hexachloro-4-[¹³ C ₁₂]biphenylol | M4H159 | 5 µg/ml |
| 2,2',3,3',4',5,5'-Heptachloro-4-[¹³ C ₁₂]biphenylol | M4H172 | 5 µg/ml |
| 2,2',3,4',5,5',6-Heptachloro-4-[¹³ C ₁₂]biphenylol | M4H187 | 5 µg/ml |



MHPCB-MXA; HRGC/LRMS TIC Chromatogram

NATIVE METHOXY-CHLOROBIPHENYLS (MeO-PCBs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-------------------|
| 4PM79 | 3,3',4',5-Tetrachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM97 | 2,2',3,4',5'-Pentachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM101 | 2,2',4,5,5'-Pentachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M107 | 2,3,3',4',5-Pentachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM108 | 2,3,3',4,5'-Pentachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 3M118 | 2,3',4,4',5-Pentachloro-3-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM120 | 2,3',4,5,5'-Pentachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM127 | 3,3',4,5,5'-Pentachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM130 | 2,2',3,3',4',5-Hexachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M134 | 2,2',3,3',5,6-Hexachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 3PM138 | 2,2',3',4,4',5-Hexachloro-3-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M146 | 2,2',3,4',5,5'-Hexachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 33PDM155 | 2,2',4,4',6,6'-Hexachloro-3,3'-dimethoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM159 | 2,3,3',4,5,5'-Hexachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M162 | 2,3,3',4',5,5'-Hexachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M163 | 2,3,3',4',5,6-Hexachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM172 | 2,2',3,3',4,5,5'-Heptachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M177 | 2,2',3,3',4',5,6-Heptachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M178 | 2,2',3,3',5,5',6-Heptachloro-4-methoxybiphenyl | 1.2 ml 48.5 µg/ml |
| 3PM180 | 2,2',3,4,4',5,5'-Heptachloro-3'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 3PM182 | 2,2',3,4,4',5,6'-Heptachloro-3'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 3PM183 | 2,2',3',4,4',5,6'-Heptachloro-3-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 3PM184 | 2,2',3,4,4',6,6'-Heptachloro-3'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M187 | 2,2',3,4',5,5',6-Heptachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M193 | 2,3,3',4',5,5',6-Heptachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM198 | 2,2',3,3',4,5,5',6-Octachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM199 | 2,2',3,3',4',5,5',6-Octachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM200 | 2,2',3,3',4,5,6,6'-Octachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM201 | 2,2',3,3',4',5,6,6'-Octachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4M202 | 2,2',3,3',5,5',6,6'-Octachloro-4-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 44PDM202 | 2,2',3,3',5,5',6,6'-Octachloro-4,4'-dimethoxybiphenyl | 1.2 ml 50 µg/ml |
| 3PM203 | 2,2',3,4,4',5,5',6-Octachloro-3'-methoxybiphenyl | 1.2 ml 50 µg/ml |
| 4PM208 | 2,2',3,3',4,5,5',6,6'-Nonachloro-4'-methoxybiphenyl | 1.2 ml 50 µg/ml |

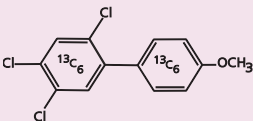
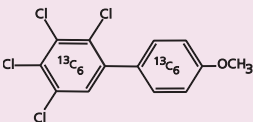
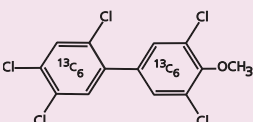
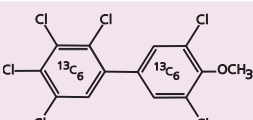
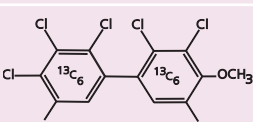
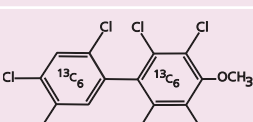
NATIVE METHOXY-CHLOROBIPHENYLS: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---|---|------------|
| MPCB-MXA | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 2,3,4,5-Tetrachloro-4'-methoxybiphenyl | 4PM61 | 5 µg/ml |
| 2,3',4,5,5'-Pentachloro-4'-methoxybiphenyl | 4PM120 | 5 µg/ml |
| 2,2',4,4',6,6'-Hexachloro-3,3'-dimethoxybiphenyl | 33PDM155 | 5 µg/ml |
| 2,2',3,4,4',6,6'-Heptachloro-3'-methoxybiphenyl | 3PM184 | 5 µg/ml |
| 2,2',3,3',5,5',6,6'-Octachloro-4-methoxybiphenyl | 4M202 | 5 µg/ml |
| MPCB-MXB | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 3,3',4',5-Tetrachloro-4-methoxybiphenyl | 4PM79 | 5 µg/ml |
| 2,2',4,5,5'-Pentachloro-4'-methoxybiphenyl | 4PM101 | 5 µg/ml |
| 2,2',3,3',5,6-Hexachloro-4-methoxybiphenyl | 4M134 | 5 µg/ml |
| 2,2',3,3',5,5',6-Heptachloro-4-methoxybiphenyl | 4M178 | 4.85 µg/ml |
| 2,2',3,3',4',5,6,6'-Octachloro-4-methoxybiphenyl | 4PM201 | 5 µg/ml |
| MPCB-MXC | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 2,3,4,4',5-Pentachloro-2'-methoxybiphenyl | 2PM114 | 5 µg/ml |
| 2,2',3,4',5,5'-Hexachloro-4-methoxybiphenyl | 4M146 | 5 µg/ml |
| 2,2',3,4,4',5,6'-Heptachloro-3'-methoxybiphenyl | 3PM182 | 5 µg/ml |
| 2,2',3,4,4',5,5',6-Octachloro-3'-methoxybiphenyl | 3PM203 | 5 µg/ml |
| 2,2',3,3',4,5,5',6,6'-Nonachloro-4'-methoxybiphenyl | 4PM208 | 5 µg/ml |
| MPCB-MXD | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 2,3',4,4',5-Pentachloro-3-methoxybiphenyl | 3M118 | 5 µg/ml |
| 2,2',3',4,4',5-Hexachloro-3-methoxybiphenyl | 3PM138 | 5 µg/ml |
| 2,2',3',4,4',5,6'-Heptachloro-3-methoxybiphenyl | 3PM183 | 5 µg/ml |
| 2,2',3,3',4,5,5',6-Octachloro-4'-methoxybiphenyl | 4PM198 | 5 µg/ml |
| MPCB-MXE | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 2,3,3',4,5'-Pentachloro-4'-methoxybiphenyl | 4PM108 | 5 µg/ml |
| 2,2',3,3',4',5-Hexachloro-4-methoxybiphenyl | 4PM130 | 5 µg/ml |
| 2,2',3,4',5,5',6-Heptachloro-4-methoxybiphenyl | 4M187 | 5 µg/ml |
| 2,2',3,3',4',5,5',6-Octachloro-4-methoxybiphenyl | 4PM199 | 5 µg/ml |
| MPCB-MXF | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| 2,3,3',4',5-Pentachloro-4-methoxybiphenyl | 4M107 | 5 µg/ml |
| 2,3,3',4',5,6-Hexachloro-4-methoxybiphenyl | 4M163 | 5 µg/ml |
| 2,2',3,3',4',5,6-Heptachloro-4-methoxybiphenyl | 4M177 | 5 µg/ml |
| 2,2',3,3',4,5,6,6'-Octachloro-4'-methoxybiphenyl | 4PM200 | 5 µg/ml |

NATIVE METHOXY-CHLOROBIPHENYLS: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|----------|
| MPCB-MXG | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| | 2,2',3,4',5'-Pentachloro-4'-methoxybiphenyl | 4PM97 |
| | 2,3,3',4,5,5'-Hexachloro-4'-methoxybiphenyl | 4PM159 |
| | 2,2',3,4,4',5,5'-Heptachloro-3'-methoxybiphenyl | 3PM180 |
| | 2,2',3,3',5,5',6,6'-Octachloro-4,4'-dimethoxybiphenyl | 44PDM202 |
| MPCB-MXH | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| | 3,3',4,5,5'-Pentachloro-4'-methoxybiphenyl | 4PM127 |
| | 2,2',3,3',4,5,5'-Heptachloro-4'-methoxybiphenyl | 4PM172 |
| MPCB-MXI | Native Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| | 2,3,3',4',5,5'-Hexachloro-4'-methoxybiphenyl | 4M162 |
| | 2,3,3',4',5,5',6-Heptachloro-4'-methoxybiphenyl | 4M193 |

MASS-LABELLED METHOXY-CHLOROBIPHENYLS

| Catalogue Number | Product |
|------------------|---|
| M4M29 |  <p>2,4,5-Trichloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M4M61 |  <p>2,3,4,5-Tetrachloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M4M120 |  <p>2,3',4,5,5'-Pentachloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M4M159 |  <p>2,3,3',4,5,5'-Hexachloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M4M172 |  <p>2,2',3,3',4,5,5'-Heptachloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |
| M4M187 |  <p>2,2',3,4',5,5',6-Heptachloro-4'-methoxy[¹³C₁₂]biphenyl 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene</p> |

* Unless stated otherwise, Isotopic Purities of these compounds are 99% or greater.

MASS-LABELLED METHOXY-CHLOROBIPHENYLS: SOLUTION/MIXTURE

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|--|----------|
| MMPCB-MXA | Mass-Labelled Chlorinated Methoxybiphenyl Solution/Mixture | 1.2 ml |
| | 2,4,5-Trichloro-4'-methoxy[¹³ C ₁₂]biphenyl | M4M29 |
| | 2,3,4,5-Tetrachloro-4'-methoxy[¹³ C ₁₂]biphenyl | M4M61 |
| | 2,3',4,5,5'-Pentachloro-4'-methoxy[¹³ C ₁₂]biphenyl | M4M120 |
| | 2,3,3',4,5,5'-Hexachloro-4'-methoxy[¹³ C ₁₂]biphenyl | M4M159 |
| | 2,2',3,3',4,5,5'-Heptachloro-4'-methoxy[¹³ C ₁₂]biphenyl | M4M172 |
| | 2,2',3,4',5,5',6-Heptachloro-4-methoxy[¹³ C ₁₂]biphenyl | M4M187 |

NATIVE CHLORINATED BIPHENYLENES (PCBPs)

Chlorinated Biphenylenes have been detected in samples taken in the aftermath of PCB fires. It is thought that they may be formed as the result of incomplete combustion of the PCBs.

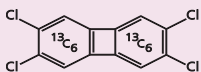
In microsomal enzyme studies, the 2,3,6,7-tetrachlorobiphenylene has shown similar potency to that of 2,3,7,8-tetrachlorodibenzo-p-dioxin in its toxicological effects.

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--------------------------------|-----------------|
| CBP-2 | 2-Chlorobiphenylene | 1.2 ml 50 µg/ml |
| CBP-23 | 2,3-Dichlorobiphenylene | 1.2 ml 50 µg/ml |
| CBP-236 | 2,3,6-Trichlorobiphenylene | 1.2 ml 50 µg/ml |
| CBP-2367 | 2,3,6,7-Tetrachlorobiphenylene | 1.2 ml 50 µg/ml |

NATIVE CHLORINATED BIPHENYLENES FOR TOXICOLOGICAL STUDIES

| Catalogue Number | Product (DMSO solution) | Qty/Conc |
|-------------------|--------------------------------|-------------------------------|
| CBP-2-D | 2-Chlorobiphenylene | 1.2 ml 1 x 10 ⁻⁴ M |
| CBP-23-D | 2,3-Dichlorobiphenylene | 1.2 ml 1 x 10 ⁻⁴ M |
| CBP-236-D | 2,3,6-Trichlorobiphenylene | 1.2 ml 1 x 10 ⁻⁴ M |
| CBP-2367-D | 2,3,6,7-Tetrachlorobiphenylene | 1.2 ml 1 x 10 ⁻⁴ M |

MASS-LABELLED CHLORINATED BIPHENYLENE

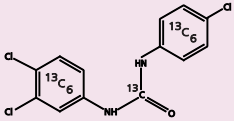
| Catalogue Number | Product |
|-------------------|---|
| MPCBP-2367 | <div style="display: flex; align-items: center;">  <div style="margin-left: 20px;"> <p>2,3,6,7-Tetrachloro[¹³C₁₂]biphenylene</p> <p>1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic Purity 99% or greater</p> </div> </div> |

NATIVE TRICLOCARBAN

Triclocarban, like Triclosan, is an antimicrobial additive used in a variety of personal care products.

| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|--|-----------------|
| TCC | N-(4-chlorophenyl)-N'-(3,4-dichlorophenyl)urea | 1.2 ml 50 µg/ml |

MASS-LABELLED TRICLOCARBAN

| Catalogue Number | Product |
|------------------|---|
| MTCC |  <p>N-(4-chloro¹³C₆]phenyl)-N'-(3,4-dichloro¹³C₆]phenyl)¹³C]urea 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; Isotopic Purity 99% or greater</p> |

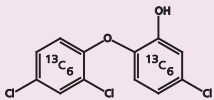
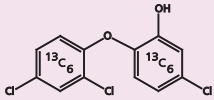
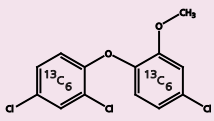
NATIVE TRICLOSAN AND METHYL TRICLOSAN

Triclosan is a widely used antibacterial and antifungal agent that has been incorporated into many common consumer products including toothpastes, deodorants, antibacterial soaps, and detergents.

The increasing use of these products over the last 30 years has led to Triclosan, and its biotransformation product, Methyl Triclosan, being found in the environment.

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------|
| TCS | 5-Chloro-2-(2,4-dichlorophenoxy)phenol | 1.2 ml 50 µg/ml |
| TCS-M | 5-Chloro-2-(2,4-dichlorophenoxy)phenol (in methanol) | 1.2 ml 50 µg/ml |
| MeTCS | 5-Chloro-2-(2,4-dichlorophenoxy)anisole | 1.2 ml 50 µg/ml |

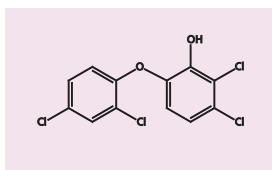
MASS-LABELLED TRICLOSAN AND METHYL TRICLOSAN

| Catalogue Number | Product |
|------------------|---|
| MTCS |  <p>5-Chloro-2-(2,4-dichloro¹³C₆]phenoxy)¹³C₆]phenol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic Purity 99% or greater</p> |
| MTCS-M |  <p>5-Chloro-2-(2,4-dichloro¹³C₆]phenoxy)¹³C₆]phenol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol; Isotopic Purity 99% or greater</p> |
| MMeTCS |  <p>5-Chloro-2-(2,4-dichloro¹³C₆]phenoxy)[1,2,3,4,5,6-¹³C₆]anisole 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane; Isotopic Purity 99% or greater</p> |

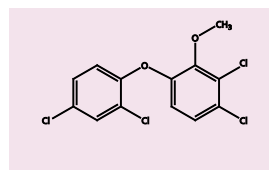
CHLORINATED DERIVATIVES OF TRICLOSAN AND METHYL TRICLOSAN

The reaction of Triclosan with free chlorine in water, or during wastewater treatment processes, may result in the formation of further chlorinated Triclosan isomers.

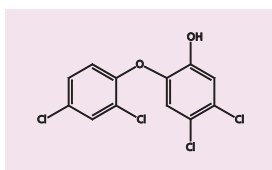
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-----------------|
| 6TCS | 2,3-Dichloro-6-(2,4-dichlorophenoxy)phenol (6-Chlorotriclosan) | 1.2 ml 50 µg/ml |
| 6MeTCS | 2,3-Dichloro-6-(2,4-dichlorophenoxy)anisole (6-Chloro-methyltriclosan) | 1.2 ml 50 µg/ml |
| 4TCS | 4,5-Dichloro-2-(2,4-dichlorophenoxy)phenol (4-Chlorotriclosan) | 1.2 ml 50 µg/ml |
| 4MeTCS | 4,5-Dichloro-2-(2,4-dichlorophenoxy)anisole (4-Chloro-methyltriclosan) | 1.2 ml 50 µg/ml |
| 46TCS | 2,3,4-Trichloro-6-(2,4-dichlorophenoxy)phenol (4,6-Dichlorotriclosan) | 1.2 ml 50 µg/ml |
| 46MeTCS | 2,3,4-Trichloro-6-(2,4-dichlorophenoxy)anisole (4,6-Dichloro-methyltriclosan) | 1.2 ml 50 µg/ml |



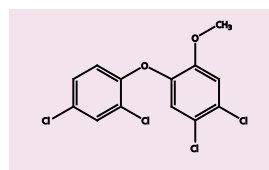
2,3-Dichloro-6-(2,4-dichlorophenoxy)phenol



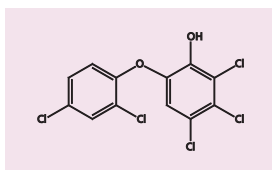
2,3-Dichloro-6-(2,4-dichlorophenoxy)anisole



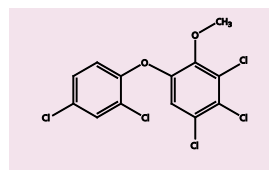
4,5-Dichloro-2-(2,4-dichlorophenoxy)phenol



4,5-Dichloro-2-(2,4-dichlorophenoxy)anisole



2,3,4-Trichloro-6-(2,4-dichlorophenoxy)phenol

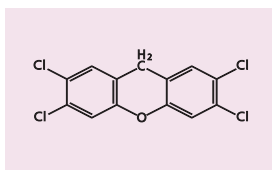


2,3,4-Trichloro-6-(2,4-dichlorophenoxy)anisole

NATIVE CHLOROANTHENE

The tetrachloroxanthene was originally suspected to be an environmental contaminant due to pulp bleaching.

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|-----------------------------|-----------------|
| XE-2367-S | 2,3,6,7-Tetrachloroxanthene | 1.2 ml 50 µg/ml |



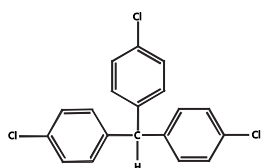
2,3,6,7-Tetrachloroxanthene

TRIS(4-CHLOROPHENYL)METHANE AND TRIS(4-CHLOROPHENYL)METHANOL; NATIVE AND MASS-LABELLED

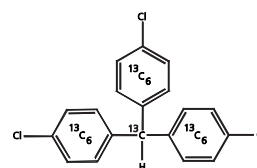
These two compounds have been found in a variety of environmental samples including fish, marine mammals and birds.

It is possible that they may originate from DDT or other agrochemicals.

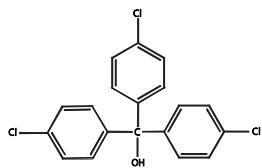
| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|--|-----------------|
| T4CPM | Tris(4-chlorophenyl)methane | 1.2 ml 50 µg/ml |
| T4CPME | Tris(4-chlorophenyl)methanol | 1.2 ml 50 µg/ml |
| MT4CPM | Tris(4-chlorophenyl)methane- ¹³ C ₁₉ ; 99 atom% ¹³ C ₁₉ | 1.2 ml 50 µg/ml |
| MT4CPME | Tris(4-chlorophenyl)methanol- ¹³ C ₁₉ ; 99 atom% ¹³ C ₁₉ | 1.2 ml 50 µg/ml |



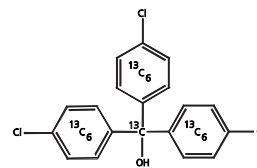
Tris(4-chlorophenyl)methane



Tris(4-chlorophenyl)methane-¹³C₁₉



Tris(4-chlorophenyl)methanol



Tris(4-chlorophenyl)methanol-¹³C₁₉

NATIVE CHLORINATED NAPHTHALENES (PCNs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---|-----------------|
| PN-2S | 2-Chloronaphthalene | 1.2 ml 50 µg/ml |
| PN-6S | 1,5-Dichloronaphthalene | 1.2 ml 50 µg/ml |
| PN-12S | 2,7-Dichloronaphthalene | 1.2 ml 50 µg/ml |
| PN-13S | 1,2,3-Trichloronaphthalene | 1.2 ml 50 µg/ml |
| PN-27S | 1,2,3,4-Tetrachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-28S | 1,2,3,5-Tetrachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-31S | 1,2,3,8-Tetrachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-46S | 1,4,5,8-Tetrachloronaphthalene (96% pure) | 1.2 ml 48 µg/ml |
| PN-48S | 2,3,6,7-Tetrachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-50S | 1,2,3,4,6-Pentachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-52S | 1,2,3,5,7-Pentachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-53S | 1,2,3,5,8-Pentachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-66S | 1,2,3,4,6,7-Hexachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-69S | 1,2,3,5,7,8-Hexachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-72S | 1,2,4,5,7,8-Hexachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-73S | 1,2,3,4,5,6,7-Heptachloronaphthalene | 1.2 ml 50 µg/ml |
| PN-75S | Octachloronaphthalene | 1.2 ml 50 µg/ml |

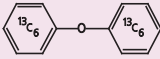
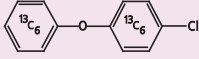
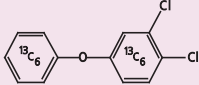
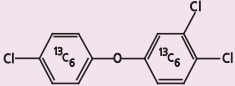
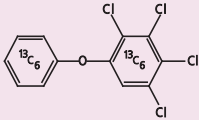
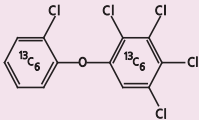
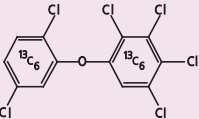
NATIVE CHLORINATED NAPHTHALENES: SOLUTION/MIXTURES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------------------------|-----------------------------|-----------|
| PCN-MXA | Native PCN Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 2-Chloronaphthalene | 2 | 5 µg/ml |
| 1,5-Dichloronaphthalene | 6 | 5 µg/ml |
| 1,2,3-Trichloronaphthalene | 13 | 5 µg/ml |
| 1,2,3,5-Tetrachloronaphthalene | 28 | 5 µg/ml |
| 1,2,3,5,7-Pentachloronaphthalene | 52 | 5 µg/ml |
| 1,2,3,4,6,7-Hexachloronaphthalene | 66 | 5 µg/ml |
| 1,2,3,4,5,6,7-Heptachloronaphthalene | 73 | 5 µg/ml |
| Octachloronaphthalene | 75 | 5 µg/ml |
| PCN-MXC | Native PCN Solution/Mixture | 1.2 ml |
| | IUPAC | |
| 1,2,3,4-Tetrachloronaphthalene | 27 | 5 µg/ml |
| 1,2,5,6-Tetrachloronaphthalene | 36 | 4.5 µg/ml |
| 1,4,5,8-Tetrachloronaphthalene | 46 | 4.8 µg/ml |
| 2,3,6,7-Tetrachloronaphthalene | 48 | 4.9 µg/ml |
| 1,2,3,4,6-Pentachloronaphthalene | 50 | 5 µg/ml |
| 1,2,3,5,8-Pentachloronaphthalene | 53 | 5 µg/ml |
| 1,2,3,5,7,8-Hexachloronaphthalene | 69 | 5 µg/ml |
| 1,2,4,5,7,8-Hexachloronaphthalene | 72 | 5 µg/ml |

NATIVE CHLORINATED DIPHENYL ETHERS (PCDEs)

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------------------|-----------------|
| DPE-0 | Diphenyl ether | 1.2 ml 50 µg/ml |
| DPE-3 | 4-Chlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-15 | 4,4'-Dichlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-28 | 2,4,4'-Trichlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-74 | 2,4,4',5-Tetrachlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-77 | 3,3',4,4'-Tetrachlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-99 | 2,2',4,4',5-Pentachlorodiphenyl ether | 1.2 ml 50 µg/ml |
| DPE-209 | Decachlorodiphenyl ether | 1.2 ml 50 µg/ml |

MASS-LABELLED CHLORINATED DIPHENYL ETHERS

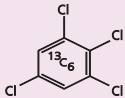
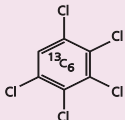
| Catalogue Number | Product |
|------------------|---|
| MCDE-0 |  <p>[¹³C₁₂]Diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-3 |  <p>4-Chloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-12 |  <p>3,4-Dichloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-37 |  <p>3,4,4'-Trichloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-61 |  <p>2,3,4,5-Tetrachloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-86 |  <p>2,2',3,4,5-Pentachloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-141 |  <p>2,2',3,4,5,5'-Hexachloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCDE-180 |  <p>2,2',3,4,4',5,5'-Heptachloro[¹³C₁₂]diphenyl ether 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, Isotopic Purities of these compounds are 99% or greater.

NATIVE CHLOROBENZENES: SOLUTION/MIXTURE

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|----------------------------|---------------------------------------|-----------|
| CBS | Native Chlorobenzene Solution/Mixture | 1.2 ml |
| Chlorobenzene | | 100 µg/ml |
| 1,2-Dichlorobenzene | | 100 µg/ml |
| 1,3-Dichlorobenzene | | 100 µg/ml |
| 1,4-Dichlorobenzene | | 100 µg/ml |
| 1,2,3-Trichlorobenzene | | 100 µg/ml |
| 1,2,4-Trichlorobenzene | | 100 µg/ml |
| 1,3,5-Trichlorobenzene | | 100 µg/ml |
| 1,2,3,4-Tetrachlorobenzene | | 100 µg/ml |
| 1,2,3,5-Tetrachlorobenzene | | 100 µg/ml |
| 1,2,4,5-Tetrachlorobenzene | | 100 µg/ml |
| Pentachlorobenzene | | 100 µg/ml |
| Hexachlorobenzene | | 100 µg/ml |

MASS-LABELLED CHLOROBENZENES

| Catalogue Number | Product |
|-------------------|---|
| MBZ-1235 |  1,2,3,5-Tetrachloro[¹³ C ₆]benzene 1.2 ml; 100 µg/ml (±5.0 µg/ml); in isooctane |
| MCBZ-12345 |  1,2,3,4,5-Pentachloro[¹³ C ₆]benzene 1.2 ml; 100 µg/ml (±5.0 µg/ml); in isooctane |

* Unless stated otherwise, Isotopic Purities of these compounds are 99% or greater.

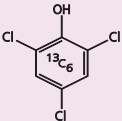
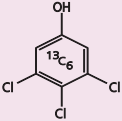
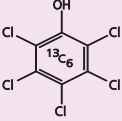
MASS-LABELLED CHLOROBENZENES: SOLUTION/MIXTURE

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|--|--|-----------|
| MCBS | Mass-Labelled Chlorobenzene Solution/Mixture | 1.2 ml |
| Chloro[¹³ C ₆]benzene | | 100 µg/ml |
| 1,4-Dichloro[¹³ C ₆]benzene | | 100 µg/ml |
| 1,2,3-Trichloro[¹³ C ₆]benzene | | 100 µg/ml |
| 1,2,3,4-Tetrachloro[¹³ C ₆]benzene | | 100 µg/ml |
| Pentachloro[¹³ C ₆]benzene | | 100 µg/ml |
| Hexachloro[¹³ C ₆]benzene | | 100 µg/ml |

NATIVE CHLOROPHENOLS: SOLUTION/MIXTURE

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|---------------------------|--------------------------------------|-----------|
| CPS | Native Chlorophenol Solution/Mixture | 1.2 ml |
| 2-Chlorophenol | | 100 µg/ml |
| 3-Chlorophenol | | 100 µg/ml |
| 4-Chlorophenol | | 100 µg/ml |
| 2,3-Dichlorophenol | | 100 µg/ml |
| 2,4-Dichlorophenol | | 100 µg/ml |
| 2,5-Dichlorophenol | | 100 µg/ml |
| 2,6-Dichlorophenol | | 100 µg/ml |
| 3,4-Dichlorophenol | | 100 µg/ml |
| 3,5-Dichlorophenol | | 100 µg/ml |
| 2,3,4-Trichlorophenol | | 100 µg/ml |
| 2,3,5-Trichlorophenol | | 100 µg/ml |
| 2,3,6-Trichlorophenol | | 100 µg/ml |
| 2,4,5-Trichlorophenol | | 100 µg/ml |
| 2,4,6-Trichlorophenol | | 100 µg/ml |
| 3,4,5-Trichlorophenol | | 100 µg/ml |
| 2,3,4,5-Tetrachlorophenol | | 100 µg/ml |
| 2,3,4,6-Tetrachlorophenol | | 100 µg/ml |
| 2,3,5,6-Tetrachlorophenol | | 100 µg/ml |
| Pentachlorophenol | | 100 µg/ml |

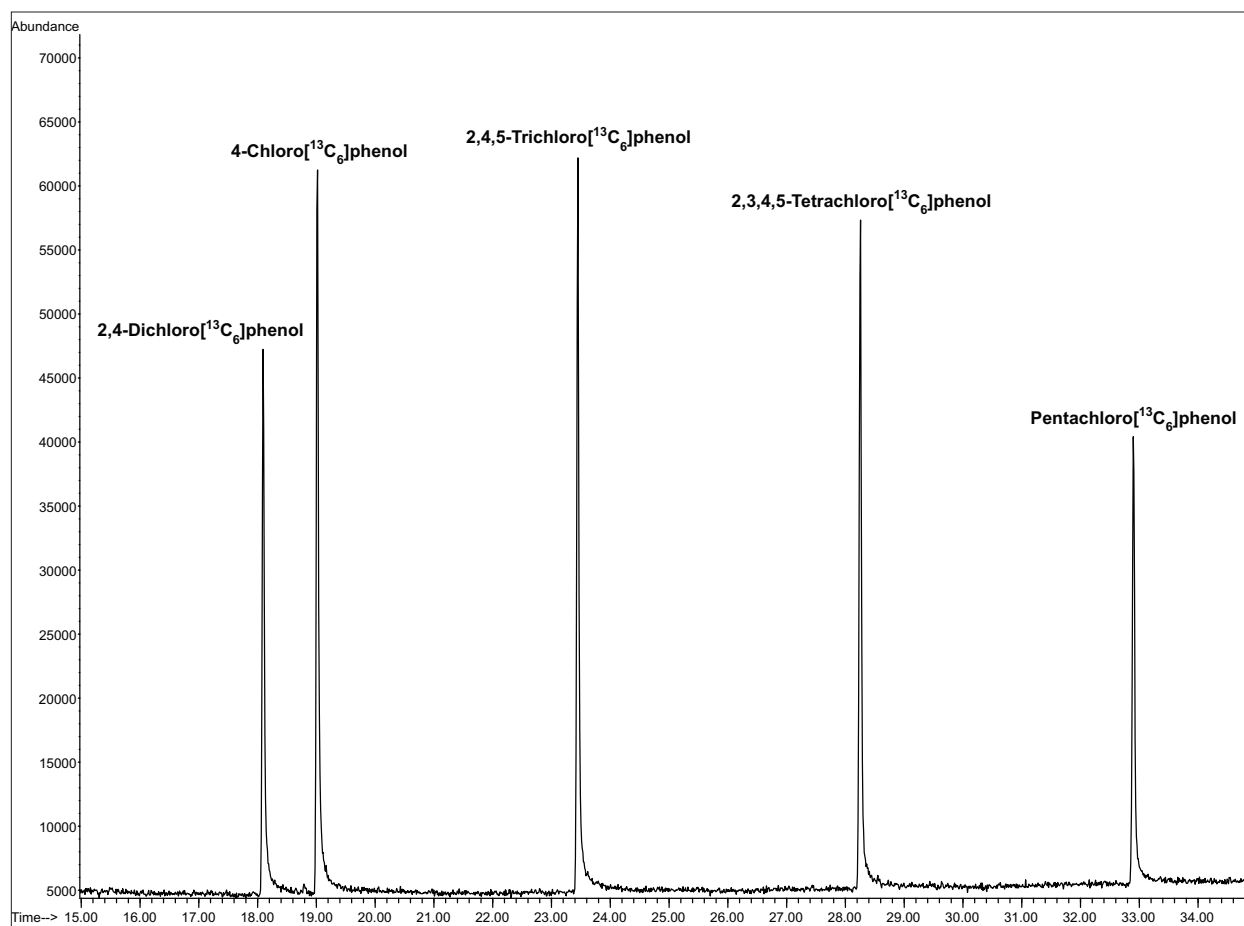
MASS-LABELLED CHLOROPHENOLS

| Catalogue Number | Product |
|------------------|---|
| MCP-246 |  <p>2,4,6-Trichloro[¹³C₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in isooctane</p> |
| MCP-345 |  <p>3,4,5-Trichloro[¹³C₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in isooctane</p> |
| MCP-23456 |  <p>Pentachloro[¹³C₆]phenol 1.2 ml; 100 µg/ml (±5.0 µg/ml); in isooctane</p> |

* Unless stated otherwise, Isotopic Purities of these compounds are 99% or greater.

MASS-LABELLED CHLOROPHENOLS: SOLUTION/MIXTURE

| Catalogue Number | Product (isooctane solution) | Qty/Conc |
|---|---|-----------|
| MCPS | Mass-Labelled Chlorophenol Solution/Mixture | 1.2 ml |
| 4-Chloro[¹³ C ₆]phenol | | 100 µg/ml |
| 2,4-Dichloro[¹³ C ₆]phenol | | 100 µg/ml |
| 2,4,5-Trichloro[¹³ C ₆]phenol | | 100 µg/ml |
| 2,3,4,5-Tetrachloro[¹³ C ₆]phenol | | 100 µg/ml |
| Pentachloro[¹³ C ₆]phenol | | 100 µg/ml |



MCPS; HRMS/LRMS TIC Chromatogram (30m DB-5 column: 0.25 mm id, 0.25 µm film thickness).

NATIVE MELAMINE AND CYANURIC ACID

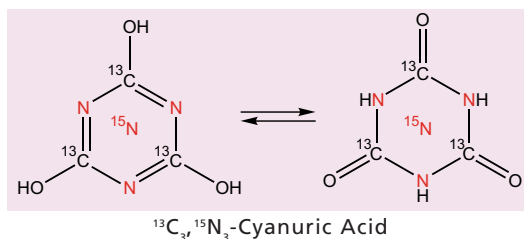
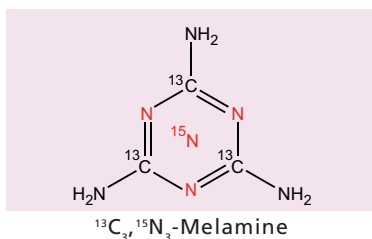
| Catalogue Number | Product (water solution) | Qty/Conc |
|------------------|--------------------------|------------------|
| MEL | Melamine | 1.2 ml 100 µg/ml |
| CYA | Cyanuric Acid | 1.2 ml 100 µg/ml |

MASS-LABELLED MELAMINE

| Catalogue Number | Product (water solution) | Qty/Conc |
|------------------|--|------------------|
| M3-MEL | $^{13}\text{C}_3$ -Melamine | 1.2 ml 100 µg/ml |
| M6-MEL | $^{13}\text{C}_3, ^{15}\text{N}_3$ -Melamine | 1.2 ml 100 µg/ml |

MASS-LABELLED CYANURIC ACID

| Catalogue Number | Product (water solution) | Qty/Conc |
|------------------|---|------------------|
| M3-CYA | $^{13}\text{C}_3$ -Cyanuric Acid | 1.2 ml 100 µg/ml |
| M6-CYA | $^{13}\text{C}_3, ^{15}\text{N}_3$ -Cyanuric Acid | 1.2 ml 100 µg/ml |

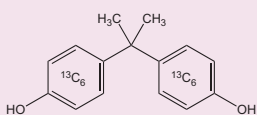


NATIVE BISPHENOL-A

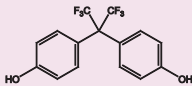
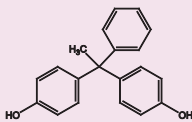
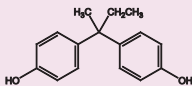
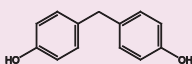
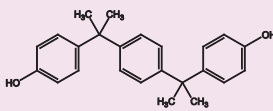
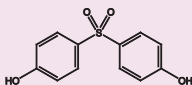
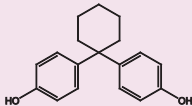
| Catalogue Number | Product (methanol solution) | Qty/Conc |
|------------------|---------------------------------|-----------------|
| BPA | 2,2-Bis(4-hydroxyphenyl)propane | 1.2 ml 50 µg/ml |

MASS-LABELLED BISPHENOL-A

| Catalogue Number | Product |
|------------------|--|
| MBPA | 2,2-Bis(4-hydroxy- $^{13}\text{C}_6$ -phenyl)propane 1.2 ml; 50 µg/ml (± 2.5 µg/ml); in methanol; Isotopic Purity 99% or greater [$^{13}\text{C}_{12}$ -rings] |



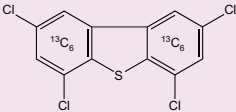
NATIVE BISPHENOL ANALOGUES

| Catalogue Number | Product |
|------------------|--|
| BPAF |  <p>2,2-Bis(4-hydroxyphenyl)hexafluoropropane 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPAP |  <p>4,4'-(1-Phenylethylidene)bisphenol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPB |  <p>2,2-Bis(4-hydroxyphenyl)butane 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPF |  <p>4,4'-Dihydroxydiphenylmethane 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPP |  <p>4,4'-(1,4-Phenylenediisopropylidene)bisphenol 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPS |  <p>Bis(4-hydroxyphenyl) sulfone 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |
| BPZ |  <p>1,1-Bis(4-hydroxyphenyl)cyclohexane 1.2 ml; 50 µg/ml (±2.5 µg/ml); in methanol</p> |

NATIVE TETRACHLORODIBENZOTHIOPHENES

| Catalogue Number | Product (toluene solution) | Qty/Conc |
|------------------|-------------------------------------|-----------------|
| TCDT-83 | 2,3,7,8-Tetrachlorodibenzothiophene | 1.2 ml 50 µg/ml |
| TCDT-85 | 2,4,6,8-Tetrachlorodibenzothiophene | 1.2 ml 50 µg/ml |

MASS-LABELLED TETRACHLORODIBENZOTHIOPHENE

| Catalogue Number | Product |
|------------------|---|
| MTCDT-85 |  <p>2,4,6,8-Tetrachloro[¹³C₁₂]dibenzothiophene 1.2 ml; 50 µg/ml (±2.5 µg/ml); in toluene; Isotopic Purity 99% or greater</p> |

NATIVE CHLORINATED CARBAZOLES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|----------------------------------|-----------------|
| CCZ-3 | 3-Chloro-9H-carbazole | 1.2 ml 50 µg/ml |
| CCZ-36 | 3,6-Dichloro-9H-carbazole | 1.2 ml 50 µg/ml |
| CCZ-1368 | 1,3,6,8-Tetrachloro-9H-carbazole | 1.2 ml 50 µg/ml |
| CCZ-2367 | 2,3,6,7-Tetrachloro-9H-carbazole | 1.2 ml 50 µg/ml |

NATIVE BROMINATED CARBAZOLES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|------------------|---------------------------------|-----------------|
| BCZ-3 | 3-Bromo-9H-carbazole | 1.2 ml 50 µg/ml |
| BCZ-27 | 2,7-Dibromo-9H-carbazole | 1.2 ml 50 µg/ml |
| BCZ-36 | 3,6-Dibromo-9H-carbazole | 1.2 ml 50 µg/ml |
| BCZ-136 | 1,3,6-Tribromo-9H-carbazole | 1.2 ml 50 µg/ml |
| BCZ-1368 | 1,3,6,8-Tetrabromo-9H-carbazole | 1.2 ml 50 µg/ml |

NATIVE BROMO/CHLORO CARBAZOLES

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|--------------------|---------------------------------------|-----------------|
| 1-B-36-CCZ | 1-Bromo-3,6-dichloro-9H-carbazole | 1.2 ml 50 µg/ml |
| 18-B-36-CCZ | 1,8-Dibromo-3,6-dichloro-9H-carbazole | 1.2 ml 50 µg/ml |

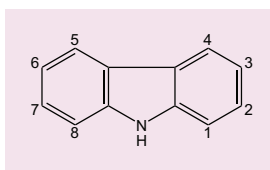
NATIVE HALOGENATED CARBAZOLES: SOLUTION/MIXTURE

| Catalogue Number | Product (nonane solution) | Qty/Conc |
|---------------------------------------|--|-----------|
| CBCZ-MXB | Native Halogenated Carbazoles Solution/Mixture | 1.2 ml |
| 3-Chloro-9H-carbazole | | 2.5 µg/ml |
| 3,6-Dichloro-9H-carbazole | | 2.5 µg/ml |
| 1,3,6,8-Tetrachloro-9H-carbazole | | 2.5 µg/ml |
| 2,3,6,7-Tetrachloro-9H-carbazole | | 2.5 µg/ml |
| 3-Bromo-9H-carbazole | | 2.5 µg/ml |
| 2,7-Dibromo-9H-carbazole | | 2.5 µg/ml |
| 3,6-Dibromo-9H-carbazole | | 2.5 µg/ml |
| 1,3,6-Tribromo-9H-carbazole | | 2.5 µg/ml |
| 1,3,6,8-Tetrabromo-9H-carbazole | | 2.5 µg/ml |
| 1-Bromo-3,6-dichloro-9H-carbazole | | 2.5 µg/ml |
| 1,8-Dibromo-3,6-dichloro-9H-carbazole | | 2.5 µg/ml |

MASS-LABELLED CHLORINATED CARBAZOLES

| Catalogue Number | Product |
|------------------|---|
| MCCZ-36 |  <p>3,6-Dichloro-9H-[¹³C₁₂]carbazole 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |
| MCCZ-1368 |  <p>1,3,6,8-Tetrachloro-9H-[¹³C₁₂]carbazole 1.2 ml; 50 µg/ml (±2.5 µg/ml); in nonane</p> |

* Unless stated otherwise, Isotopic Purities of these compounds are 99% or greater.



General Structure and Numbering System for Carbazoles



Ingrid O'Gorman
South River, Ontario, Canada

APPENDIX

Some of the information provided in this Appendix can also be found in our **Reference and Handling Guide: GC/MS Characterization and Analysis of Selected Halogenated Aromatic Compounds**, which is available separately in hard copy and is also posted on our website. On the following pages you will find:

Guidelines for the Use and Handling of Wellington Products

General Structure and Numbering System of Selected Aromatic Hydrocarbons

Number of Possible Isomers for Selected Halogenated Aromatic Compounds

Molecular Weights for Selected Chlorinated and Brominated Aromatic Hydrocarbons

Exact Mass & Relative Ion Abundance of Selected Chlorinated Aromatic Hydrocarbons

Molecular Ion Clusters for Chlorinated Aromatic Hydrocarbons & Brominated Aromatic Hydrocarbons

Exact Mass & Relative Ion Abundance of Selected Brominated Aromatic Hydrocarbons

Systematic Numbering of Chlorinated Dibenzo-p-dioxins, Chlorinated Dibenzofurans, Chlorinated Biphenyls, and Chlorinated Naphthalenes

We have also prepared a **Reference and Handling Guide for Perfluoroalkyl Compounds**, which can be downloaded from our website, and a more concise **Quick Reference Guide for Perfluoroalkyl Compounds**, which is available in hard copy.

If you would like to receive copies of either Reference Guide, please contact us or one of our distributors. In addition, if you have any ideas on how we can improve these Reference Guides, or if you have suggestions for future guides, please contact us.



GUIDELINES FOR THE USE AND HANDLING OF WELLINGTON PRODUCTS

HAZARDS

The majority of our products are halogenated aromatic hydrocarbons in solution in organic solvents such as nonane, toluene and isooctane. Although the maximum concentration is 100 µg/ml, that is 0.01% (w/v), these compounds must be considered toxic and potentially carcinogenic and should be handled accordingly.

With all of our products due care should be exercised to prevent human contact and ingestion. The absence of a toxicity warning for any of our products must not be interpreted as an indication that there is no possible health hazard.



NOTE:

THESE MATERIALS SHOULD ONLY BE USED BY PERSONNEL TRAINED IN THE HANDLING OF HAZARDOUS CHEMICALS.

ALL PROCEDURES SHOULD BE PERFORMED IN A FUME HOOD AND SUITABLE GLOVES, EYE PROTECTION AND CLOTHING SHOULD BE WORN AT ALL TIMES.

RECEIPT, INSPECTION, HANDLING AND STORAGE

Unless crystalline material is provided, all of our reference standard solutions come in flame-sealed, pre-scored amber glass ampoules. Upon receipt, inspect the ampoules for breakage and leakage and then store them upright until needed. The ampoules can be stored at ambient temperature until opened unless other storage requirements are stated on the Certificate of Analysis.

Prior to opening, allow the solution to drain into the bottom of the ampoule, lightly tapping the ampoule if necessary. Using the plastic ampoule collar provided, hold the ampoule upright and snap the top off, breaking away from the body.

Transfer the solution to an amber glass container that can be tightly sealed for storage. To prevent evaporation of the solvent, it is suggested that this solution, and subsequent mixtures and/or dilutions, be stored at refrigerator temperatures.

GUIDELINES FOR THE USE AND HANDLING OF WELLINGTON PRODUCTS

DISPOSAL

All waste materials generated during the use of these solutions should be treated as hazardous in accordance with national and regional regulations. A licensed disposal company should be employed. Some options for the destruction of these materials include high temperature incineration, photolysis, or chemical treatment using reagents such as sodium naphthalene or KPEG reagent. Literature references for some of these methods can be provided upon request.

ACCURACY

Each of our stock solutions is prepared from crystalline material that has been well characterized as to its structure and purity.

The crystalline material is weighed using microbalances that are externally calibrated using NIST-traceable weights. Solutions are prepared by completely dissolving the crystalline material in ultrapure, distilled-in-glass solvents. The volumetric flasks used for this purpose, and the pipets used for subsequent preparation of dilutions and mixtures, are all of class A tolerance and NIST-traceable.

The maximum percent relative combined uncertainty for solution preparation is calculated to be $\pm 5\%$.

INTERLABORATORY CERTIFICATION

Wellington continues to submit its standards for independent interlaboratory testing and certification. Since 1991, our standards have been tested in 30 international round-robins.

To date, solutions of the compounds listed below have been repeatedly tested and the approximate total number of analyses are given.

- 2,3,7,8 - substituted PCDDs and PCDFs1500 HRMS analyses
- Dioxin-like (WHO) PCB congeners1000 HRMS analyses
- PBDEs200 HRMS analyses
- PFCs200 LCMS analyses

The overall averages of the data received for all of the compounds were found to be well within $\pm 10\%$ of the design values.

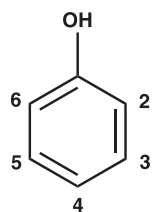
EXPIRY DATE/SHELF LIFE

In order to accurately determine the shelf life of products such as ours, testing must reveal significant degradation or loss in concentration of the particular analyte. In comparing freshly prepared solutions to older solutions by GC/MS or LC/MS, we have not detected any significant changes. Many of these older solutions were prepared and ampouled more than 15 years ago. Thus our stability studies, as they should, remain ongoing.

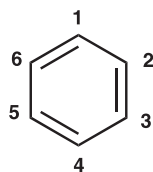
For our products where the expiry date on the C of A states, "stability studies ongoing", we consider that our reference standard solutions retain their accuracy for a period of 5 years from delivery in the unopened ampoule.

NOTE: The predominant degradation pathway for our compounds is likely photolysis and thus protection from light is critical.

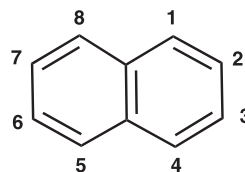
GENERAL STRUCTURE AND NUMBERING SYSTEM OF SELECTED AROMATIC HYDROCARBONS



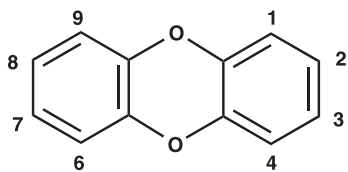
phenol



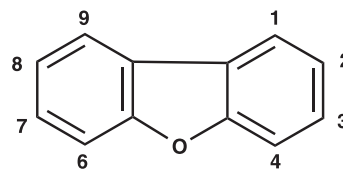
benzene



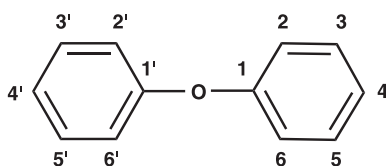
naphthalene



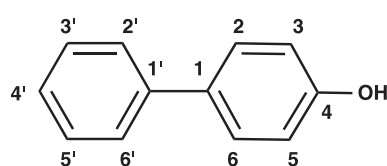
dibenzo-p-dioxin



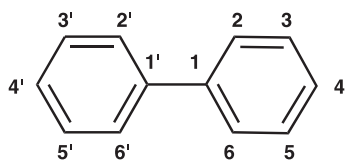
dibenzofuran



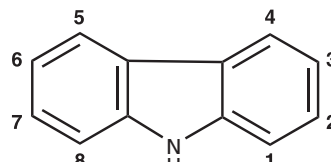
diphenyl ether



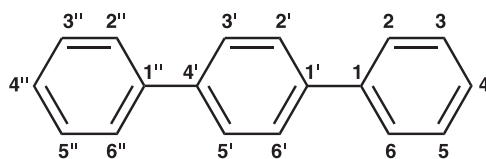
4-hydroxybiphenyl



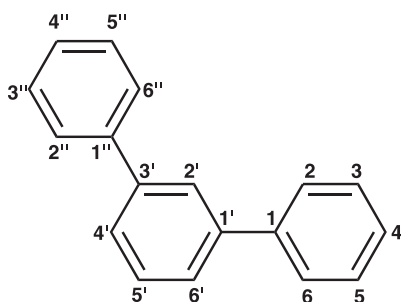
biphenyl



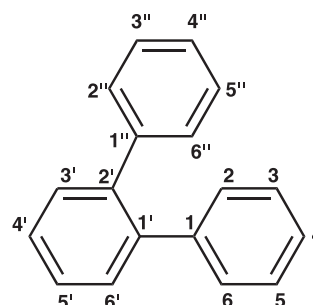
carbazole



p-terphenyl



m-terphenyl



o-terphenyl

NUMBER OF POSSIBLE ISOMERS FOR SELECTED HALOGENATED AROMATIC COMPOUNDS

| # of X | Terphenyl | | | Biphenyl | Biphenylol | Dibenzo-p-dioxin | Dibenzofuran | Naphthalene | Benzene | Phenol |
|--------|-----------|------|------|----------|------------|------------------|--------------|-------------|---------|--------|
| | ortho | meta | para | | | | | | | |
| 1 | 5 | 6 | 4 | 3 | 19 | 2 | 4 | 2 | 1 | 3 |
| 2 | 28 | 28 | 21 | 12 | 64 | 10 | 16 | 10 | 3 | 6 |
| 3 | 80 | 87 | 55 | 24 | 136 | 14 | 28 | 14 | 3 | 6 |
| 4 | 211 | 211 | 139 | 42 | 198 | 22 | 38 | 22 | 3 | 3 |
| 5 | 355 | 382 | 226 | 46 | 198 | 14 | 28 | 14 | 1 | 1 |
| 6 | 544 | 544 | 351 | 42 | 136 | 10 | 16 | 10 | 1 | |
| 7 | 596 | 638 | 358 | 24 | 64 | 2 | 4 | 2 | | |
| 8 | 544 | 544 | 351 | 12 | 19 | 1 | 1 | 1 | | |
| 9 | 355 | 382 | 226 | 3 | 3 | | | | | |
| 10 | 211 | 211 | 139 | 1 | | | | | | |
| 11 | 80 | 87 | 55 | | | | | | | |
| 12 | 28 | 28 | 21 | | | | | | | |
| 13 | 5 | 6 | 4 | | | | | | | |
| 14 | 1 | 1 | 1 | | | | | | | |

X= Halogen (does not apply to mixed halogenated compounds)
For diphenyl ethers use the biphenyl values

MOLECULAR WEIGHTS FOR SELECTED CHLORINATED AND BROMINATED AROMATIC HYDROCARBONS

| # of Cl/Br | PCTs | PCBs | PCDEs | PCDDs | PCDFs | PCNs | CBs | CPs | PBBs | PBDEs | PBDDs | PBDFs |
|------------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|--------|
| 0 | 230.31 | 154.21 | 170.21 | 184.19 | 168.19 | 128.17 | 78.11 | 94.11 | 154.21 | 170.21 | 184.19 | 168.19 |
| 1 | 264.75 | 188.66 | 204.66 | 218.64 | 202.64 | 162.62 | 112.56 | 128.56 | 233.11 | 249.11 | 263.09 | 247.09 |
| 2 | 299.20 | 223.10 | 239.10 | 253.08 | 237.09 | 197.06 | 147.00 | 163.00 | 312.00 | 328.00 | 341.99 | 325.99 |
| 3 | 333.64 | 257.55 | 273.55 | 287.53 | 271.53 | 231.51 | 181.45 | 197.45 | 390.90 | 406.90 | 420.88 | 404.88 |
| 4 | 368.09 | 291.99 | 307.99 | 321.97 | 305.98 | 265.95 | 215.89 | 231.89 | 469.80 | 485.80 | 499.78 | 483.78 |
| 5 | 402.53 | 326.44 | 342.44 | 356.42 | 340.42 | 300.40 | 250.34 | 266.34 | 548.69 | 564.69 | 578.67 | 562.68 |
| 6 | 436.98 | 360.88 | 376.88 | 390.86 | 374.87 | 334.84 | 284.78 | | 627.59 | 643.59 | 657.57 | 641.57 |
| 7 | 471.42 | 395.33 | 411.33 | 425.31 | 409.31 | 369.29 | | | 706.48 | 722.48 | 736.47 | 720.47 |
| 8 | 505.87 | 429.77 | 445.77 | 459.75 | 443.76 | 403.73 | | | 785.38 | 801.38 | 815.36 | 799.36 |
| 9 | 540.31 | 464.22 | 480.22 | | | | | | 864.28 | 880.28 | | |
| 10 | 574.76 | 498.66 | 514.66 | | | | | | 943.17 | 959.17 | | |
| 11 | 609.20 | | | | | | | | | | | |
| 12 | 643.65 | | | | | | | | | | | |
| 13 | 678.09 | | | | | | | | | | | |
| 14 | 712.54 | | | | | | | | | | | |

Note: The molecular weight for PCHBs ($C_{12}H_{9-n}Cl_nOH$) is the same as the PCDEs ($C_{12}H_{10-n}Cl_nO$), but the maximum # of Chlorines is one less for the PCHBs.

PCTs = polychlorinated terphenyls, PCBs = polychlorinated biphenyls, PCDEs = polychlorinated diphenyl ethers
PCHBs = polychlorinated hydroxybiphenyls (biphenylols), PCDDs = polychlorinated dibenzo-p-dioxins,
PCDFs = polychlorinated dibenzofurans, PCNs = polychlorinated naphthalenes, CBs = chlorobenzenes, CPs = chlorophenols,
PBBs = polybrominated biphenyls, PBDEs = polybrominated diphenyl ethers,
PBDDs = polybrominated dibenzo-p-dioxins, PBDFs = polybrominated dibenzofurans

EXACT MASS & RELATIVE ION ABUNDANCE OF SELECTED CHLORINATED AROMATIC HYDROCARBONS

| # of Cl | PCTs | | | PCBs | | | PCHBs | | | PCDDs | | |
|---------|-------------------------------|---|-------------------------------|-------------------------------|---|-------------------------------|-------------------------------|---|-------------------------------|-------------------------------|---|-------------------------------|
| | ¹² C ₁₈ | C ₁₈ H _{14-n} Cl _n | ¹³ C ₁₈ | ¹² C ₁₂ | C ₁₂ H _{10-n} Cl _n | ¹³ C ₁₂ | ¹² C ₁₂ | C ₁₂ H _{9-n} Cl _n OH | ¹³ C ₁₂ | ¹² C ₁₂ | C ₁₂ H _{8-n} Cl _n O ₂ | ¹³ C ₁₂ |
| | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass |
| 0 | 230.1096 | 100 | 248.1699 | 154.0783 | 100 | 166.1185 | 170.0732 | 100 | 182.1134 | 184.0524 | 100 | 196.0927 |
| 1 | 264.0706 | 100 | 282.1310 | 188.0393 | 100 | 200.0795 | 204.0342 | 100 | 216.0745 | 218.0135 | 100 | 230.0537 |
| | 266.0676 | 34.4 | 284.1280 | 190.0363 | 33.2 | 202.0766 | 206.0312 | 33.5 | 218.0715 | 220.0105 | 33.7 | 232.0508 |
| 2 | 298.0316 | 100 | 316.0920 | 222.0003 | 100 | 234.0406 | 237.9952 | 100 | 250.0355 | 251.9745 | 100 | 264.0147 |
| | 300.0287 | 66.8 | 318.0890 | 223.9974 | 65.6 | 236.0376 | 239.9923 | 65.9 | 252.0325 | 253.9715 | 66.1 | 266.0118 |
| | 302.0257 | 11.8 | 320.0861 | 225.9944 | 11.0 | 238.0347 | 241.9893 | 11.2 | 254.0296 | 255.9686 | 11.3 | 268.0088 |
| 3 | 331.9926 | 100 | 350.0530 | 255.9613 | 100 | 268.0016 | 271.9562 | 100 | 283.9965 | 285.9355 | 100 | 297.9758 |
| | 333.9897 | 99.2 | 352.0501 | 257.9584 | 98.0 | 269.9986 | 273.9533 | 98.2 | 285.9936 | 287.9326 | 98.5 | 299.9728 |
| | 335.9867 | 33.4 | 354.0471 | 259.9554 | 32.3 | 271.9957 | 275.9503 | 32.5 | 287.9906 | 289.9296 | 32.7 | 301.9699 |
| | 337.9838 | 4.0 | 356.0442 | 261.9525 | 3.7 | 273.9927 | 277.9474 | 3.7 | 289.9877 | 291.9267 | 3.8 | 303.9669 |
| 4 | 365.9537 | 76.0 | 384.0141 | 289.9224 | 76.7 | 301.9626 | 305.9173 | 76.5 | 317.9575 | 319.8965 | 76.4 | 331.9368 |
| | 367.9507 | 100 | 386.0111 | 291.9194 | 100 | 303.9597 | 307.9143 | 100 | 319.9546 | 321.8936 | 100 | 333.9339 |
| | 369.9478 | 49.8 | 388.0082 | 293.9165 | 49.1 | 305.9567 | 309.9114 | 49.3 | 321.9516 | 323.8906 | 49.4 | 335.9309 |
| | 371.9448 | 11.3 | 390.0052 | 295.9135 | 10.8 | 307.9538 | 311.9084 | 10.9 | 323.9487 | 325.8877 | 11.0 | 337.9280 |
| 5 | 399.9147 | 61.0 | 417.9751 | 323.8834 | 61.4 | 335.9237 | 339.8783 | 61.3 | 351.9186 | 353.8576 | 61.3 | 365.8978 |
| | 401.9117 | 100 | 419.9721 | 325.8804 | 100 | 337.9207 | 341.8754 | 100 | 353.9156 | 355.8546 | 100 | 367.8949 |
| | 403.9088 | 66.0 | 421.9692 | 327.8775 | 65.3 | 339.9178 | 343.8724 | 65.4 | 355.9127 | 357.8517 | 65.5 | 369.8919 |
| | 405.9058 | 22.0 | 423.9662 | 329.8745 | 21.4 | 341.9148 | 345.8695 | 21.5 | 357.9097 | 359.8487 | 21.6 | 371.8890 |
| 6 | 433.8757 | 50.9 | 451.9361 | 357.8444 | 51.2 | 369.8847 | 373.8393 | 51.2 | 385.8796 | 387.8186 | 51.1 | 399.8589 |
| | 435.8728 | 100 | 453.9332 | 359.8415 | 100 | 371.8817 | 375.8364 | 100 | 387.8766 | 389.8156 | 100 | 401.8559 |
| | 437.8698 | 82.1 | 455.9302 | 361.8385 | 81.5 | 373.8788 | 377.8334 | 81.6 | 389.8737 | 391.8127 | 81.7 | 403.8530 |
| | 439.8669 | 36.2 | 457.9273 | 363.8356 | 35.5 | 375.8758 | 379.8305 | 35.6 | 391.8707 | 393.8097 | 35.8 | 405.8500 |
| 7 | 467.8367 | 43.7 | 485.8971 | 391.8054 | 43.9 | 403.8457 | 407.8004 | 43.9 | 419.8406 | 421.7796 | 43.9 | 433.8199 |
| | 469.8338 | 100 | 487.8942 | 393.8025 | 100 | 405.8428 | 409.7974 | 100 | 421.8377 | 423.7767 | 100 | 435.8169 |
| | 471.8308 | 98.3 | 489.8912 | 395.7995 | 97.7 | 407.8398 | 411.7945 | 97.8 | 423.8347 | 425.7737 | 97.9 | 437.8140 |
| | 473.8279 | 53.9 | 491.8883 | 397.7966 | 53.1 | 409.8369 | 413.7915 | 53.3 | 425.8318 | 427.7708 | 53.4 | 439.8110 |
| | 475.8249 | 17.9 | 493.8853 | 399.7936 | 17.4 | 411.8339 | 415.7886 | 17.5 | 427.8288 | 429.7678 | 17.6 | 441.8081 |
| 8 | 501.7978 | 33.5 | 519.8582 | 425.7665 | 33.8 | 437.8067 | 441.7614 | 33.7 | 453.8016 | 455.7407 | 33.7 | 467.7809 |
| | 503.7948 | 87.4 | 521.8552 | 427.7635 | 87.8 | 439.8038 | 443.7584 | 87.7 | 455.7987 | 457.7377 | 87.6 | 469.7780 |
| | 505.7919 | 100 | 523.8523 | 429.7606 | 100 | 441.8008 | 445.7555 | 100 | 457.7957 | 459.7348 | 100 | 471.7750 |
| | 507.7889 | 65.6 | 525.8493 | 431.7576 | 65.2 | 443.7979 | 447.7525 | 65.2 | 459.7928 | 461.7318 | 65.3 | 473.7721 |
| | 509.7860 | 27.0 | 527.8464 | 433.7547 | 26.6 | 445.7949 | 449.7496 | 26.7 | 461.7898 | 463.7289 | 26.8 | 475.7691 |
| 9 | 535.7588 | 26.1 | 553.8192 | 459.7275 | 26.3 | 471.7678 | 475.7224 | 26.3 | 487.7627 | | | |
| | 537.7559 | 76.5 | 555.8162 | 461.7246 | 76.9 | 473.7648 | 477.7195 | 76.8 | 489.7597 | | | |
| | 539.7529 | 100 | 557.8133 | 463.7216 | 100 | 475.7619 | 479.7165 | 100 | 491.7568 | | | |
| | 541.7500 | 76.4 | 559.8103 | 465.7187 | 75.9 | 477.7589 | 481.7136 | 76.0 | 493.7538 | | | |
| | 543.7470 | 37.6 | 561.8074 | 467.7157 | 37.1 | 479.7560 | 483.7106 | 37.2 | 495.7509 | | | |
| 10 | 569.7198 | 20.9 | 587.7802 | 493.6885 | 21.1 | 505.7288 | | | | | | |
| | 571.7169 | 68.1 | 589.7773 | 495.6856 | 68.4 | 507.7258 | | | | | | |
| | 573.7139 | 100 | 591.7743 | 497.6826 | 100 | 509.7229 | | | | | | |
| | 575.7110 | 87.2 | 593.7714 | 499.6797 | 86.7 | 511.7199 | | | | | | |
| | 577.7080 | 50.0 | 595.7684 | 501.6767 | 49.4 | 513.7170 | | | | | | |
| 11 | 603.6809 | 17.1 | 621.7412 | | | | | | | | | |
| | 605.6779 | 61.3 | 623.7383 | | | | | | | | | |
| | 607.6750 | 100 | 625.7353 | | | | | | | | | |
| | 609.6720 | 98.0 | 627.7324 | | | | | | | | | |
| | 611.6691 | 64.1 | 629.7294 | | | | | | | | | |
| | 613.6661 | 29.4 | 631.7265 | | | | | | | | | |
| 12 | 637.6419 | 13.1 | 655.7023 | | | | | | | | | |
| | 639.6389 | 51.3 | 657.6993 | | | | | | | | | |
| | 641.6360 | 92.0 | 659.6964 | | | | | | | | | |
| | 643.6330 | 100 | 661.6934 | | | | | | | | | |
| | 645.6301 | 73.5 | 663.6905 | | | | | | | | | |
| | 647.6271 | 38.5 | 665.6875 | | | | | | | | | |
| 13 | 671.6029 | 10.1 | 689.6633 | | | | | | | | | |
| | 673.6000 | 42.8 | 691.6604 | | | | | | | | | |
| | 675.5970 | 83.7 | 693.6574 | | | | | | | | | |
| | 677.5941 | 100 | 695.6545 | | | | | | | | | |
| | 679.5911 | 81.6 | 697.6515 | | | | | | | | | |
| | 681.5882 | 48.0 | 699.6486 | | | | | | | | | |
| 14 | 705.5639 | 8.0 | 723.6243 | | | | | | | | | |
| | 707.5610 | 36.3 | 725.6214 | | | | | | | | | |
| | 709.5580 | 76.7 | 727.6184 | | | | | | | | | |
| | 711.5551 | 100 | 729.6155 | | | | | | | | | |
| | 713.5521 | 89.7 | 731.6125 | | | | | | | | | |
| | 715.5492 | 58.6 | 733.6096 | | | | | | | | | |

EXACT MASS & RELATIVE ION ABUNDANCE OF SELECTED CHLORINATED AROMATIC HYDROCARBONS

| # of Cl | PCDFs | | | PCNs | | | CBs | | | CPs | | |
|---------|-------------------------------|--|-------------------------------|-------------------------------|--|-------------------------------|------------------------------|---|------------------------------|------------------------------|--|------------------------------|
| | ¹² C ₁₂ | C ₁₂ H _{8-n} Cl _n O | ¹³ C ₁₂ | ¹² C ₁₀ | C ₁₀ H _{8-n} Cl _n | ¹³ C ₁₀ | ¹² C ₆ | C ₆ H _{5-n} Cl _n | ¹³ C ₆ | ¹² C ₆ | C ₆ H _{5-n} Cl _n OH | ¹³ C ₆ |
| | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass |
| 0 | 168.0575 | 100 | 180.0978 | 128.0626 | 100 | 138.0962 | 78.0470 | 100 | 84.0671 | 94.0419 | 100 | 100.0620 |
| 1 | 202.0185 | 100 | 214.0588 | 162.0236 | 100 | 172.0572 | 112.0080 | 100 | 118.0281 | 128.0029 | 100 | 134.0230 |
| | 204.0156 | 33.5 | 216.0559 | 164.0207 | 33.0 | 174.0542 | 114.0050 | 32.6 | 120.0252 | 129.9999 | 32.8 | 136.0201 |
| 2 | 235.9796 | 100 | 248.0198 | 195.9847 | 100 | 206.0182 | 145.9690 | 100 | 151.9891 | 161.9639 | 100 | 167.9841 |
| | 237.9766 | 65.8 | 250.0169 | 197.9817 | 65.4 | 208.0153 | 147.9661 | 65.0 | 153.9862 | 163.9610 | 65.2 | 169.9811 |
| | 239.9737 | 11.2 | 252.0139 | 199.9788 | 10.9 | 210.0123 | 149.9631 | 10.6 | 155.9832 | 165.9580 | 10.8 | 171.9782 |
| 3 | 269.9406 | 100 | 281.9809 | 229.9457 | 100 | 239.9792 | 179.9300 | 100 | 185.9502 | 195.9249 | 100 | 201.9451 |
| | 271.9376 | 98.2 | 283.9779 | 231.9427 | 97.8 | 241.9763 | 181.9271 | 97.4 | 187.9472 | 197.9220 | 97.6 | 203.9421 |
| | 273.9347 | 32.5 | 285.9750 | 233.9398 | 32.0 | 243.9733 | 183.9241 | 31.7 | 189.9443 | 199.9190 | 31.9 | 205.9392 |
| | 275.9317 | 3.7 | 287.9720 | 235.9368 | 3.6 | 245.9704 | 185.9212 | 3.5 | 191.9413 | 201.9161 | 3.5 | 207.9362 |
| 4 | 303.9016 | 76.5 | 315.9419 | 263.9067 | 76.8 | 273.9403 | 213.8911 | 77.1 | 219.9112 | 229.8860 | 76.9 | 235.9061 |
| | 305.8987 | 100 | 317.9389 | 265.9038 | 100 | 275.9373 | 215.8881 | 100 | 221.9082 | 231.8830 | 100 | 237.9032 |
| | 307.8957 | 49.2 | 319.9360 | 267.9008 | 49.0 | 277.9344 | 217.8852 | 48.7 | 223.9053 | 233.8801 | 48.8 | 239.9002 |
| | 309.8928 | 10.9 | 321.9330 | 269.8979 | 10.7 | 279.9314 | 219.8822 | 10.6 | 225.9023 | 235.8771 | 10.7 | 241.8973 |
| 5 | 337.8627 | 61.3 | 349.9029 | 297.8677 | 61.5 | 307.9013 | 247.8521 | 61.7 | 253.8722 | 263.8470 | 61.6 | 269.8671 |
| | 339.8597 | 100 | 351.9000 | 299.8648 | 100 | 309.8983 | 249.8491 | 100 | 255.8693 | 265.8441 | 100 | 271.8642 |
| | 341.8568 | 65.4 | 353.8970 | 301.8618 | 65.1 | 311.8954 | 251.8462 | 64.9 | 257.8663 | 267.8411 | 65.0 | 273.8612 |
| | 343.8538 | 21.5 | 355.8941 | 303.8589 | 21.3 | 313.8924 | 253.8432 | 21.1 | 259.8634 | 269.8382 | 21.2 | 275.8583 |
| 6 | 371.8237 | 51.2 | 383.8639 | 331.8288 | 51.3 | 341.8623 | 281.8131 | 51.4 | 287.8332 | | | |
| | 373.8207 | 100 | 385.8610 | 333.8258 | 100 | 343.8594 | 283.8102 | 100 | 289.8303 | | | |
| | 375.8178 | 81.6 | 387.8580 | 335.8229 | 81.3 | 345.8564 | 285.8072 | 81.1 | 291.8273 | | | |
| | 377.8148 | 35.6 | 389.8551 | 337.8199 | 35.3 | 347.8535 | 287.8043 | 35.1 | 293.8244 | | | |
| 7 | 405.7847 | 43.9 | 417.8250 | 365.7898 | 44.0 | 375.8233 | | | | | | |
| | 407.7818 | 100 | 419.8220 | 367.7868 | 100 | 377.8204 | | | | | | |
| | 409.7788 | 97.8 | 421.8191 | 369.7839 | 97.5 | 379.8174 | | | | | | |
| | 411.7759 | 53.3 | 423.8161 | 371.7809 | 52.9 | 381.8145 | | | | | | |
| | 413.7729 | 17.5 | 425.8132 | 373.7780 | 17.3 | 383.8115 | | | | | | |
| 8 | 439.7457 | 33.7 | 451.7860 | 399.7508 | 33.9 | 409.7844 | | | | | | |
| | 441.7428 | 87.7 | 453.7830 | 401.7479 | 87.9 | 411.7814 | | | | | | |
| | 443.7398 | 100 | 455.7801 | 403.7449 | 100 | 413.7785 | | | | | | |
| | 445.7369 | 65.2 | 457.7771 | 405.7420 | 65.0 | 415.7755 | | | | | | |
| | 447.7339 | 26.7 | 459.7742 | 407.7390 | 26.5 | 417.7726 | | | | | | |

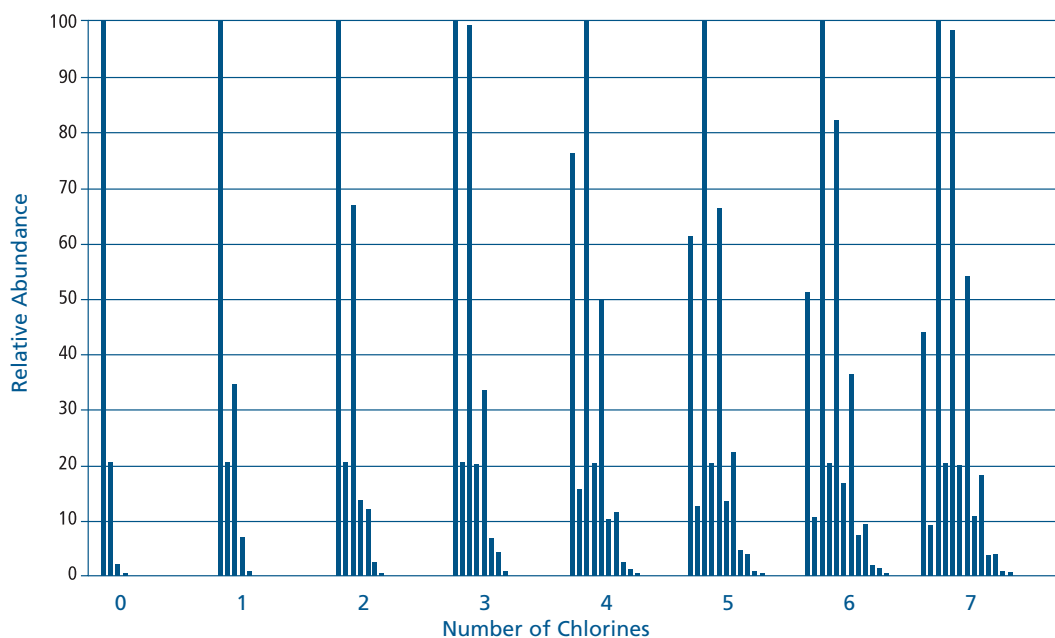
- PCTs** = polychlorinated terphenyls
PCBs = polychlorinated biphenyls
PCHBs = polychlorinated hydroxybiphenyls
PCDDs = polychlorinated dibenzo-p-dioxins
PCDFs = polychlorinated dibenzofurans
PCNs = polychlorinated naphthalenes
CBs = chlorobenzenes
CPs = chlorophenols

Accurate masses: ¹²C=12.000000, ¹³C=13.003355, ¹H=1.007825, ³⁵Cl= 34.968853, ³⁷Cl=36.965903, ¹⁶O=15.994915

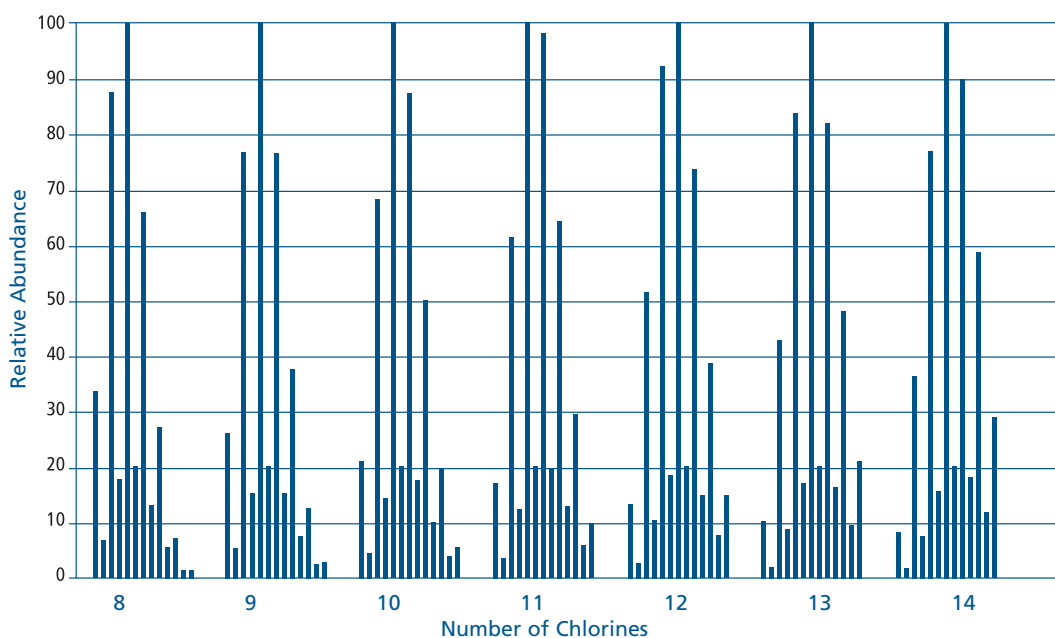
Relative abundances of isotopes were determined using the method described in: Pretsch,Clerc,Seibl,Simon, Tables of Spectral Data for Structure Determination of Organic Compounds, Springer-Verlag, 1983.

The following natural isotopic abundances were used in all calculations: ¹²C=98.89%, ¹³C=1.11%, ¹H=99.985%, ²H=0.015%, ³⁵Cl=75.53%, ³⁷Cl=24.47%, ¹⁶O=99.759%, ¹⁷O=0.037%, ¹⁸O=0.204%.

MOLECULAR ION CLUSTERS FOR CHLORINATED AROMATIC HYDROCARBONS

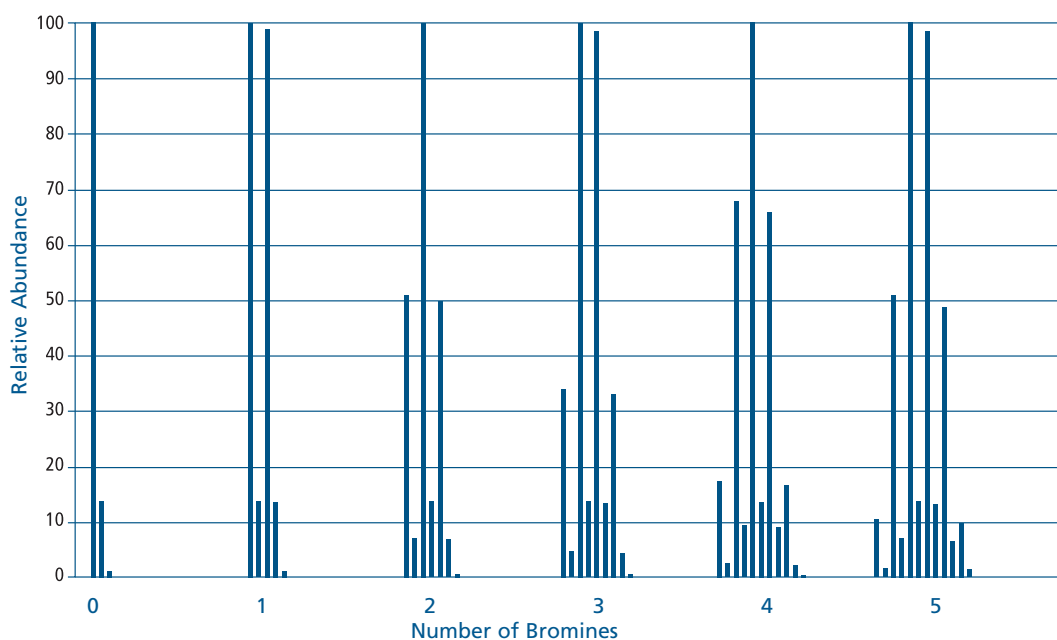


MOLECULAR ION CLUSTERS FOR CHLORINATED AROMATIC HYDROCARBONS

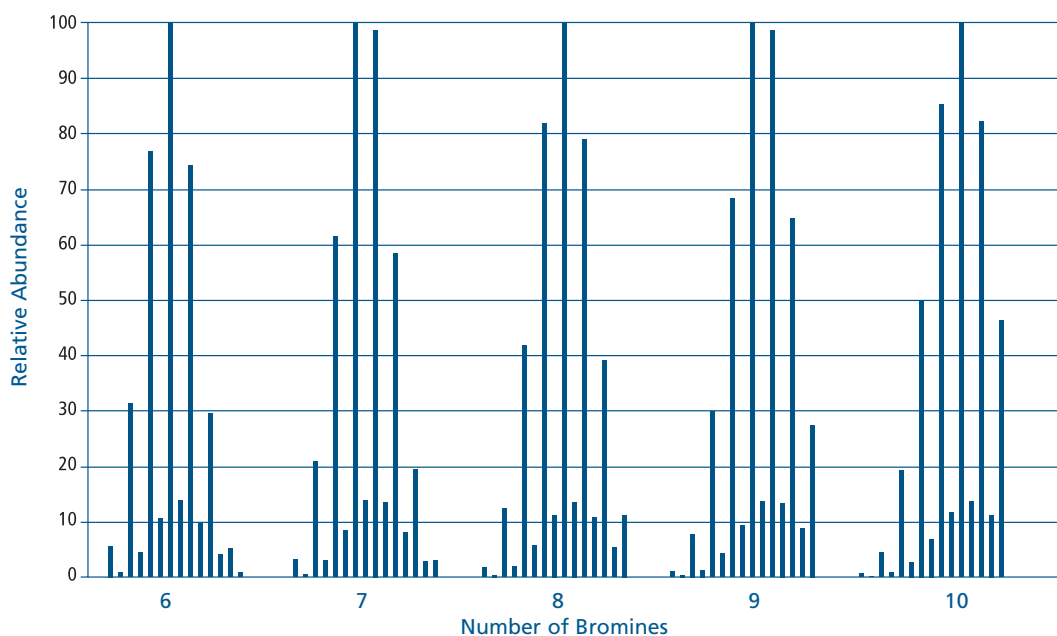


Ions shown are M, M+1, M+2, M+3, M+4, M+5, M+6, M+7, M+8, M+9, M+10, M+11, M+12, and are representative of chlorinated terphenyls ($C_{18}H_{14-n}Cl_n$)

MOLECULAR ION CLUSTERS FOR BROMINATED AROMATIC HYDROCARBONS



MOLECULAR ION CLUSTERS FOR BROMINATED AROMATIC HYDROCARBONS



Ions shown are M, M+1, M+2, M+3, M+4, M+5, M+6, M+7, M+8, M+9, M+10, M+11, M+12, M+13, M+14 and are representative of brominated diphenyl ethers ($C_{12}H_{10-n}Br_nO$)

EXACT MASS & RELATIVE ION ABUNDANCE OF SELECTED BROMINATED AROMATIC HYDROCARBONS

| # of Br | PBBs | | | PBDEs | | | PBDDs | | | PBDFs | | |
|---------|-------------------------------|---|-------------------------------|-------------------------------|---|-------------------------------|-------------------------------|---|-------------------------------|-------------------------------|--|-------------------------------|
| | ¹² C ₁₂ | C ₁₂ H _{10-n} Br _n | ¹³ C ₁₂ | ¹² C ₁₂ | C ₁₂ H _{10-n} Br _n O | ¹³ C ₁₂ | ¹² C ₁₂ | C ₁₂ H _{8-n} Br _n O ₂ | ¹³ C ₁₂ | ¹² C ₁₂ | C ₁₂ H _{8-n} Br _n O | ¹³ C ₁₂ |
| | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass | Exact Mass | Relative Abundance | Exact Mass |
| 0 | 154.0783 | 100 | 166.1185 | 170.0732 | 100 | 182.1134 | 184.0524 | 100 | 196.0927 | 168.0575 | 100 | 180.0978 |
| 1 | 231.9887 | 100 | 244.0290 | 247.9836 | 100 | 260.0239 | 261.9629 | 100 | 274.0032 | 245.9680 | 100 | 258.0083 |
| | 233.9867 | 98.7 | 246.0270 | 249.9816 | 98.9 | 262.0219 | 263.9609 | 99.1 | 276.0012 | 247.9660 | 98.9 | 260.0063 |
| 2 | 309.8992 | 50.9 | 321.9395 | 325.8941 | 50.8 | 337.9344 | 339.8734 | 50.8 | 351.9136 | 323.8785 | 50.8 | 335.9187 |
| | 311.8972 | 100 | 323.9375 | 327.8921 | 100 | 339.9324 | 341.8714 | 100 | 353.9116 | 325.8765 | 100 | 337.9167 |
| | 313.8952 | 49.6 | 325.9355 | 329.8901 | 49.7 | 341.9304 | 343.8694 | 49.9 | 355.9096 | 327.8745 | 49.7 | 339.9147 |
| 3 | 387.8097 | 34.0 | 399.8499 | 403.8046 | 33.9 | 415.8449 | 417.7839 | 33.9 | 429.8241 | 401.7889 | 33.9 | 413.8292 |
| | 389.8077 | 100 | 401.8479 | 405.8026 | 100 | 417.8429 | 419.7819 | 100 | 431.8221 | 403.7869 | 100 | 415.8272 |
| | 391.8057 | 98.4 | 403.8459 | 407.8006 | 98.6 | 419.8409 | 421.7799 | 98.7 | 433.8201 | 405.7849 | 98.6 | 417.8252 |
| | 393.8037 | 32.7 | 405.8439 | 409.7986 | 32.8 | 421.8389 | 423.7779 | 33.0 | 435.8181 | 407.7829 | 32.8 | 419.8232 |
| 4 | 465.7202 | 17.3 | 477.7604 | 481.7151 | 17.3 | 493.7553 | 495.6943 | 17.3 | 507.7346 | 479.6994 | 17.3 | 491.7397 |
| | 467.7182 | 67.9 | 479.7584 | 483.7131 | 67.8 | 495.7533 | 497.6923 | 67.8 | 509.7326 | 481.6974 | 67.8 | 493.7377 |
| | 469.7162 | 100 | 481.7564 | 485.7111 | 100 | 497.7513 | 499.6903 | 100 | 511.7306 | 483.6954 | 100 | 495.7357 |
| | 471.7142 | 65.7 | 483.7544 | 487.7091 | 65.8 | 499.7493 | 501.6883 | 65.9 | 513.7286 | 485.6934 | 65.8 | 497.7337 |
| | 473.7122 | 16.4 | 485.7524 | 489.7071 | 16.5 | 501.7473 | 503.6863 | 16.6 | 515.7266 | 487.6914 | 16.5 | 499.7317 |
| 5 | 543.6306 | 10.4 | 555.6709 | 559.6255 | 10.4 | 571.6658 | 573.6048 | 10.4 | 585.6451 | 557.6099 | 10.4 | 569.6502 |
| | 545.6286 | 51.0 | 557.6689 | 561.6235 | 50.9 | 573.6638 | 575.6028 | 50.9 | 587.6431 | 559.6079 | 50.9 | 571.6482 |
| | 547.6266 | 100 | 559.6669 | 563.6215 | 100 | 575.6618 | 577.6008 | 100 | 589.6411 | 561.6059 | 100 | 573.6462 |
| | 549.6246 | 98.3 | 561.6649 | 565.6195 | 98.4 | 577.6598 | 579.5988 | 98.5 | 591.6391 | 563.6039 | 98.4 | 575.6442 |
| | 551.6226 | 48.5 | 563.6629 | 567.6175 | 48.7 | 579.6578 | 581.5968 | 48.8 | 593.6371 | 565.6019 | 48.7 | 577.6422 |
| | 553.6206 | 9.7 | 565.6609 | 569.6155 | 9.8 | 581.6558 | 583.5948 | 9.9 | 595.6351 | 567.5999 | 9.8 | 579.6402 |
| 6 | 621.5411 | 5.3 | 633.5814 | 637.5360 | 5.3 | 649.5763 | 651.5153 | 5.3 | 663.5555 | 635.5204 | 5.3 | 647.5606 |
| | 623.5391 | 31.2 | 635.5794 | 639.5340 | 31.1 | 651.5743 | 653.5133 | 31.1 | 665.5535 | 637.5184 | 31.1 | 649.5586 |
| | 625.5371 | 76.4 | 637.5774 | 641.5320 | 76.4 | 653.5723 | 655.5113 | 76.3 | 667.5515 | 639.5164 | 76.4 | 651.5566 |
| | 627.5351 | 100 | 639.5754 | 643.5300 | 100 | 655.5703 | 657.5093 | 100 | 669.5495 | 641.5144 | 100 | 653.5546 |
| | 629.5331 | 73.8 | 641.5734 | 645.5280 | 73.9 | 657.5683 | 659.5073 | 73.9 | 671.5475 | 643.5124 | 73.9 | 655.5526 |
| | 631.5311 | 29.2 | 643.5714 | 647.5260 | 29.3 | 659.5663 | 661.5053 | 29.4 | 673.5455 | 645.5104 | 29.3 | 657.5506 |
| 7 | 699.4516 | 3.0 | 711.4918 | 715.4465 | 3.0 | 727.4868 | 729.4258 | 3.0 | 741.4660 | 713.4308 | 3.0 | 725.4711 |
| | 701.4496 | 20.8 | 713.4898 | 717.4445 | 20.8 | 729.4848 | 731.4238 | 20.8 | 743.4640 | 715.4288 | 20.8 | 727.4691 |
| | 703.4476 | 61.2 | 715.4878 | 719.4425 | 61.1 | 731.4828 | 733.4218 | 61.1 | 745.4620 | 717.4268 | 61.1 | 729.4671 |
| | 705.4456 | 100 | 717.4858 | 721.4405 | 100 | 733.4808 | 735.4198 | 100 | 747.4600 | 719.4248 | 100 | 731.4651 |
| | 707.4436 | 98.2 | 719.4838 | 723.4385 | 98.3 | 735.4788 | 737.4178 | 98.4 | 749.4580 | 721.4228 | 98.3 | 733.4631 |
| | 709.4416 | 58.0 | 721.4818 | 725.4365 | 58.1 | 737.4768 | 739.4158 | 58.2 | 751.4560 | 723.4208 | 58.1 | 735.4611 |
| 8 | 777.3621 | 1.5 | 789.4023 | 793.3570 | 1.5 | 805.3972 | 807.3362 | 1.5 | 819.3765 | 791.3413 | 1.5 | 803.3816 |
| | 779.3601 | 12.1 | 791.4003 | 795.3550 | 12.1 | 807.3952 | 809.3342 | 12.1 | 821.3745 | 793.3393 | 12.1 | 805.3796 |
| | 781.3581 | 41.6 | 793.3983 | 797.3530 | 41.5 | 809.3932 | 811.3322 | 41.5 | 823.3725 | 795.3373 | 41.5 | 807.3776 |
| | 783.3561 | 81.5 | 795.3963 | 799.3510 | 81.5 | 811.3912 | 813.3302 | 81.4 | 825.3705 | 797.3353 | 81.5 | 809.3756 |
| | 785.3541 | 100 | 797.3943 | 801.3490 | 100 | 813.3892 | 815.3282 | 100 | 827.3685 | 799.3333 | 100 | 811.3736 |
| | 787.3521 | 78.6 | 799.3923 | 803.3470 | 78.7 | 815.3872 | 817.3262 | 78.7 | 829.3665 | 801.3313 | 78.7 | 813.3716 |
| | 789.3501 | 38.7 | 801.3903 | 805.3450 | 38.8 | 817.3852 | 819.3242 | 38.9 | 831.3645 | 803.3293 | 38.8 | 815.3696 |
| 9 | 855.2725 | 0.9 | 867.3128 | 871.2674 | 0.9 | 883.3077 | | | | | | |
| | 857.2705 | 7.6 | 869.3108 | 873.2654 | 7.6 | 885.3057 | | | | | | |
| | 859.2685 | 29.7 | 871.3088 | 875.2634 | 29.7 | 887.3037 | | | | | | |
| | 861.2665 | 68.0 | 873.3068 | 877.2614 | 68.0 | 889.3017 | | | | | | |
| | 863.2645 | 100 | 875.3048 | 879.2594 | 100 | 891.2997 | | | | | | |
| | 865.2625 | 98.1 | 877.3028 | 881.2574 | 98.2 | 893.2977 | | | | | | |
| | 867.2605 | 64.3 | 879.3008 | 883.2554 | 64.4 | 895.2957 | | | | | | |
| | 869.2585 | 27.2 | 881.2988 | 885.2534 | 27.3 | 897.2937 | | | | | | |
| 10 | 933.1830 | 0.4 | 945.2233 | 949.1779 | 0.4 | 961.2182 | | | | | | |
| | 935.1810 | 4.3 | 947.2213 | 951.1759 | 4.3 | 963.2162 | | | | | | |
| | 937.1790 | 19.0 | 949.2193 | 953.1739 | 18.9 | 965.2142 | | | | | | |
| | 939.1770 | 49.5 | 951.2173 | 955.1719 | 49.5 | 967.2122 | | | | | | |
| | 941.1750 | 85.0 | 953.2153 | 957.1699 | 84.9 | 969.2102 | | | | | | |
| | 943.1730 | 100 | 955.2133 | 959.1679 | 100 | 971.2082 | | | | | | |
| | 945.1710 | 81.8 | 957.2113 | 961.1659 | 81.9 | 973.2062 | | | | | | |
| | 947.1690 | 46.0 | 959.2093 | 963.1639 | 46.0 | 975.2042 | | | | | | |
| | 949.1670 | 17.0 | 961.2073 | 965.1619 | 17.1 | 977.2022 | | | | | | |

PBBs = polybrominated biphenyls
PBDEs = polybrominated diphenyl ethers
PBDDs = polybrominated dibenzo-p-dioxins
PBDFs = polybrominated dibenzofurans

Accurate masses: ¹²C=12.000000, ¹³C=13.003355, ¹H=1.007825, ⁷⁹Br= 78.918300, ⁸¹Br=80.916300, ¹⁶O=15.994915

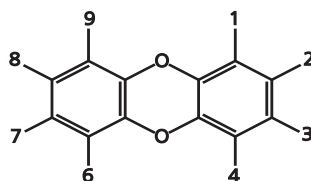
Relative abundances of isotopes were determined using the method described in: Pretsch, Clerc, Seibl, Simon, Tables of Spectral Data for Structure Determination of Organic Compounds, Springer-Verlag, 1983.

The following natural isotopic abundances were used in all calculations: ¹²C=98.89%, ¹³C=1.11%, ¹H=99.985%, ²H=0.015%, ⁷⁹Br=50.54%, ⁸¹Br=49.46%, ¹⁶O=99.759%, ¹⁷O=0.037%, ¹⁸O=0.204%.

SYSTEMATIC NUMBERING OF CHLORINATED DIBENZO-p-DIOXINS

| ID Number* | Congener | CAS Number |
|------------|-------------------------------------|------------|
| 1 | 1-Chlorodibenzo-p-dioxin | 39227-53-7 |
| 2 | 2-Chlorodibenzo-p-dioxin | 39227-54-8 |
| 3 | 1,2-Dichlorodibenzo-p-dioxin | 54536-18-4 |
| 4 | 1,3-Dichlorodibenzo-p-dioxin | 50585-39-2 |
| 5 | 1,4-Dichlorodibenzo-p-dioxin | 54536-19-5 |
| 6 | 1,6-Dichlorodibenzo-p-dioxin | 58178-38-0 |
| 7 | 1,7-Dichlorodibenzo-p-dioxin | 82291-26-7 |
| 8 | 1,8-Dichlorodibenzo-p-dioxin | 82291-27-8 |
| 9 | 1,9-Dichlorodibenzo-p-dioxin | 82291-28-9 |
| 10 | 2,3-Dichlorodibenzo-p-dioxin | 29446-15-9 |
| 11 | 2,7-Dichlorodibenzo-p-dioxin | 33857-26-0 |
| 12 | 2,8-Dichlorodibenzo-p-dioxin | 38964-22-6 |
| 13 | 1,2,3-Trichlorodibenzo-p-dioxin | 54536-17-3 |
| 14 | 1,2,4-Trichlorodibenzo-p-dioxin | 39227-58-2 |
| 15 | 1,2,6-Trichlorodibenzo-p-dioxin | 82291-29-0 |
| 16 | 1,2,7-Trichlorodibenzo-p-dioxin | 82291-30-3 |
| 17 | 1,2,8-Trichlorodibenzo-p-dioxin | 82291-31-4 |
| 18 | 1,2,9-Trichlorodibenzo-p-dioxin | 82291-32-5 |
| 19 | 1,3,6-Trichlorodibenzo-p-dioxin | 82291-33-6 |
| 20 | 1,3,7-Trichlorodibenzo-p-dioxin | 67028-17-5 |
| 21 | 1,3,8-Trichlorodibenzo-p-dioxin | 82306-61-4 |
| 22 | 1,3,9-Trichlorodibenzo-p-dioxin | 82306-62-5 |
| 23 | 1,4,6-Trichlorodibenzo-p-dioxin | 82306-63-6 |
| 24 | 1,4,7-Trichlorodibenzo-p-dioxin | 82306-64-7 |
| 25 | 1,7,8-Trichlorodibenzo-p-dioxin | 82306-65-8 |
| 26 | 2,3,7-Trichlorodibenzo-p-dioxin | 33857-28-2 |
| 27 | 1,2,3,4-Tetrachlorodibenzo-p-dioxin | 30746-58-8 |
| 28 | 1,2,3,6-Tetrachlorodibenzo-p-dioxin | 71669-25-5 |
| 29 | 1,2,3,7-Tetrachlorodibenzo-p-dioxin | 67028-18-6 |
| 30 | 1,2,3,8-Tetrachlorodibenzo-p-dioxin | 53555-02-5 |
| 31 | 1,2,3,9-Tetrachlorodibenzo-p-dioxin | 71669-26-6 |
| 32 | 1,2,4,6-Tetrachlorodibenzo-p-dioxin | 71669-27-7 |

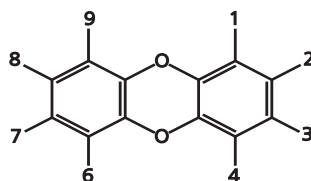
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZO-p-DIOXINS

| ID Number* | Congener | CAS Number |
|------------|--|------------|
| 33 | 1,2,4,7-Tetrachlorodibenzo-p-dioxin | 71669-28-8 |
| 34 | 1,2,4,8-Tetrachlorodibenzo-p-dioxin | 71669-29-9 |
| 35 | 1,2,4,9-Tetrachlorodibenzo-p-dioxin | 71665-99-1 |
| 36 | 1,2,6,7-Tetrachlorodibenzo-p-dioxin | 40581-90-6 |
| 37 | 1,2,6,8-Tetrachlorodibenzo-p-dioxin | 67323-56-2 |
| 38 | 1,2,6,9-Tetrachlorodibenzo-p-dioxin | 40581-91-7 |
| 39 | 1,2,7,8-Tetrachlorodibenzo-p-dioxin | 34816-53-0 |
| 40 | 1,2,7,9-Tetrachlorodibenzo-p-dioxin | 71669-23-3 |
| 41 | 1,2,8,9-Tetrachlorodibenzo-p-dioxin | 62470-54-6 |
| 42 | 1,3,6,8-Tetrachlorodibenzo-p-dioxin | 33423-92-6 |
| 43 | 1,3,6,9-Tetrachlorodibenzo-p-dioxin | 71669-24-4 |
| 44 | 1,3,7,8-Tetrachlorodibenzo-p-dioxin | 50585-46-1 |
| 45 | 1,3,7,9-Tetrachlorodibenzo-p-dioxin | 62470-53-5 |
| 46 | 1,4,6,9-Tetrachlorodibenzo-p-dioxin | 40581-93-9 |
| 47 | 1,4,7,8-Tetrachlorodibenzo-p-dioxin | 40581-94-0 |
| 48 | 2,3,7,8-Tetrachlorodibenzo-p-dioxin | 1746-01-6 |
| 49 | 1,2,3,4,6-Pentachlorodibenzo-p-dioxin | 67028-19-7 |
| 50 | 1,2,3,4,7-Pentachlorodibenzo-p-dioxin | 39227-61-7 |
| 51 | 1,2,3,6,7-Pentachlorodibenzo-p-dioxin | 71925-15-0 |
| 52 | 1,2,3,6,8-Pentachlorodibenzo-p-dioxin | 71925-16-1 |
| 53 | 1,2,3,6,9-Pentachlorodibenzo-p-dioxin | 82291-34-7 |
| 54 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin | 40321-76-4 |
| 55 | 1,2,3,7,9-Pentachlorodibenzo-p-dioxin | 71925-17-2 |
| 56 | 1,2,3,8,9-Pentachlorodibenzo-p-dioxin | 71925-18-3 |
| 57 | 1,2,4,6,7-Pentachlorodibenzo-p-dioxin | 82291-35-8 |
| 58 | 1,2,4,6,8-Pentachlorodibenzo-p-dioxin | 71998-76-0 |
| 59 | 1,2,4,6,9-Pentachlorodibenzo-p-dioxin | 82291-36-9 |
| 60 | 1,2,4,7,8-Pentachlorodibenzo-p-dioxin | 58802-08-7 |
| 61 | 1,2,4,7,9-Pentachlorodibenzo-p-dioxin | 82291-37-0 |
| 62 | 1,2,4,8,9-Pentachlorodibenzo-p-dioxin | 82291-38-1 |
| 63 | 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin | 58200-66-1 |
| 64 | 1,2,3,4,6,8-Hexachlorodibenzo-p-dioxin | 58200-67-2 |
| 65 | 1,2,3,4,6,9-Hexachlorodibenzo-p-dioxin | 58200-68-3 |

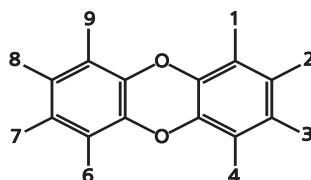
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZO-p-DIOXINS

| ID Number* | Congener | CAS Number |
|------------|---|------------|
| 66 | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin | 39227-28-6 |
| 67 | 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin | 57653-85-7 |
| 68 | 1,2,3,6,7,9-Hexachlorodibenzo-p-dioxin | 64461-98-9 |
| 69 | 1,2,3,6,8,9-Hexachlorodibenzo-p-dioxin | 58200-69-4 |
| 70 | 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin | 19408-74-3 |
| 71 | 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin | 39227-62-8 |
| 72 | 1,2,4,6,8,9-Hexachlorodibenzo-p-dioxin | 58802-09-8 |
| 73 | 1,2,3,4,6,7,8-Heptachlorodibenzo-p-dioxin | 35822-46-9 |
| 74 | 1,2,3,4,6,7,9-Heptachlorodibenzo-p-dioxin | 58200-70-7 |
| 75 | Octachlorodibenzo-p-dioxin | 3268-87-9 |

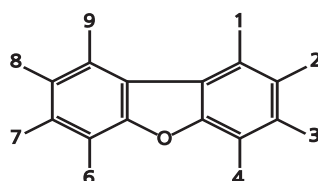
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZOFURANS

| ID Number* | Congener | CAS Number |
|------------|--------------------------|------------|
| 1 | 1-Chlorodibenzofuran | 84761-86-4 |
| 2 | 2-Chlorodibenzofuran | 51230-49-0 |
| 3 | 3-Chlorodibenzofuran | 25074-67-3 |
| 4 | 4-Chlorodibenzofuran | 74992-96-4 |
| 5 | 1,2-Dichlorodibenzofuran | 64126-85-8 |
| 6 | 1,3-Dichlorodibenzofuran | 94538-00-8 |
| 7 | 1,4-Dichlorodibenzofuran | 94538-01-9 |
| 8 | 1,6-Dichlorodibenzofuran | 74992-97-5 |
| 9 | 1,7-Dichlorodibenzofuran | 94538-02-0 |
| 10 | 1,8-Dichlorodibenzofuran | 81638-37-1 |
| 11 | 1,9-Dichlorodibenzofuran | 70648-14-5 |

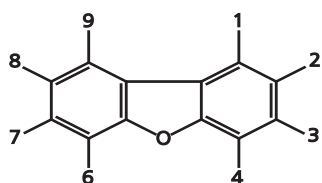
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZOFURANS

| ID Number* | Congener | CAS Number |
|------------|-----------------------------|------------|
| 12 | 2,3-Dichlorodibenzofuran | 64126-86-9 |
| 13 | 2,4-Dichlorodibenzofuran | |
| 14 | 2,6-Dichlorodibenzofuran | 60390-27-4 |
| 15 | 2,7-Dichlorodibenzofuran | 74992-98-6 |
| 16 | 2,8-Dichlorodibenzofuran | 5409-83-6 |
| 17 | 3,4-Dichlorodibenzofuran | 94570-83-9 |
| 18 | 3,6-Dichlorodibenzofuran | 74918-40-4 |
| 19 | 3,7-Dichlorodibenzofuran | 58802-21-4 |
| 20 | 4,6-Dichlorodibenzofuran | |
| 21 | 1,2,3-Trichlorodibenzofuran | 83636-47-9 |
| 22 | 1,2,4-Trichlorodibenzofuran | 24478-73-7 |
| 23 | 1,2,6-Trichlorodibenzofuran | 64560-15-2 |
| 24 | 1,2,7-Trichlorodibenzofuran | 83704-37-4 |
| 25 | 1,2,8-Trichlorodibenzofuran | 83704-34-1 |
| 26 | 1,2,9-Trichlorodibenzofuran | 83704-38-5 |
| 27 | 1,3,4-Trichlorodibenzofuran | 82911-61-3 |
| 28 | 1,3,6-Trichlorodibenzofuran | 83704-39-6 |
| 29 | 1,3,7-Trichlorodibenzofuran | 64560-16-3 |
| 30 | 1,3,8-Trichlorodibenzofuran | 76621-12-0 |
| 31 | 1,3,9-Trichlorodibenzofuran | 83704-40-9 |
| 32 | 1,4,6-Trichlorodibenzofuran | 82911-60-2 |
| 33 | 1,4,7-Trichlorodibenzofuran | 83704-41-0 |
| 34 | 1,4,8-Trichlorodibenzofuran | 64560-14-1 |
| 35 | 1,4,9-Trichlorodibenzofuran | 70648-13-4 |
| 36 | 1,6,7-Trichlorodibenzofuran | 83704-46-5 |
| 37 | 1,6,8-Trichlorodibenzofuran | 82911-59-9 |
| 38 | 1,7,8-Trichlorodibenzofuran | 58802-18-9 |
| 39 | 2,3,4-Trichlorodibenzofuran | 57117-34-7 |
| 40 | 2,3,6-Trichlorodibenzofuran | 57117-33-6 |
| 41 | 2,3,7-Trichlorodibenzofuran | 58802-17-8 |
| 42 | 2,3,8-Trichlorodibenzofuran | 57117-32-5 |
| 43 | 2,4,6-Trichlorodibenzofuran | 58802-14-5 |
| 44 | 2,4,7-Trichlorodibenzofuran | 83704-42-1 |
| 45 | 2,4,8-Trichlorodibenzofuran | 54589-71-8 |

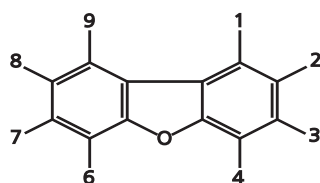
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZOFURANS

| ID Number* | Congener | CAS Number |
|------------|---------------------------------|------------|
| 46 | 2,6,7-Trichlorodibenzofuran | 83704-45-4 |
| 47 | 3,4,6-Trichlorodibenzofuran | 83704-43-2 |
| 48 | 3,4,7-Trichlorodibenzofuran | 83704-44-3 |
| 49 | 1,2,3,4-Tetrachlorodibenzofuran | 24478-72-6 |
| 50 | 1,2,3,6-Tetrachlorodibenzofuran | 83704-21-6 |
| 51 | 1,2,3,7-Tetrachlorodibenzofuran | 83704-22-7 |
| 52 | 1,2,3,8-Tetrachlorodibenzofuran | 62615-08-1 |
| 53 | 1,2,3,9-Tetrachlorodibenzofuran | 83704-23-8 |
| 54 | 1,2,4,6-Tetrachlorodibenzofuran | 71998-73-7 |
| 55 | 1,2,4,7-Tetrachlorodibenzofuran | 83719-40-8 |
| 56 | 1,2,4,8-Tetrachlorodibenzofuran | 64126-87-0 |
| 57 | 1,2,4,9-Tetrachlorodibenzofuran | 83704-24-9 |
| 58 | 1,2,6,7-Tetrachlorodibenzofuran | 83704-25-0 |
| 59 | 1,2,6,8-Tetrachlorodibenzofuran | 83710-07-0 |
| 60 | 1,2,6,9-Tetrachlorodibenzofuran | 70648-18-9 |
| 61 | 1,2,7,8-Tetrachlorodibenzofuran | 58802-20-3 |
| 62 | 1,2,7,9-Tetrachlorodibenzofuran | 83704-26-1 |
| 63 | 1,2,8,9-Tetrachlorodibenzofuran | 70648-22-5 |
| 64 | 1,3,4,6-Tetrachlorodibenzofuran | 83704-27-2 |
| 65 | 1,3,4,7-Tetrachlorodibenzofuran | 70648-16-7 |
| 66 | 1,3,4,8-Tetrachlorodibenzofuran | 92341-04-3 |
| 67 | 1,3,4,9-Tetrachlorodibenzofuran | 83704-28-3 |
| 68 | 1,3,6,7-Tetrachlorodibenzofuran | 57117-36-9 |
| 69 | 1,3,6,8-Tetrachlorodibenzofuran | 71998-72-6 |
| 70 | 1,3,6,9-Tetrachlorodibenzofuran | 83690-98-6 |
| 71 | 1,3,7,8-Tetrachlorodibenzofuran | 57117-35-8 |
| 72 | 1,3,7,9-Tetrachlorodibenzofuran | 64560-17-4 |
| 73 | 1,4,6,7-Tetrachlorodibenzofuran | 66794-59-0 |
| 74 | 1,4,6,8-Tetrachlorodibenzofuran | 82911-58-8 |
| 75 | 1,4,6,9-Tetrachlorodibenzofuran | 70648-19-0 |
| 76 | 1,4,7,8-Tetrachlorodibenzofuran | 83704-29-4 |
| 77 | 1,6,7,8-Tetrachlorodibenzofuran | 83704-33-0 |
| 78 | 2,3,4,6-Tetrachlorodibenzofuran | 83704-30-7 |
| 79 | 2,3,4,7-Tetrachlorodibenzofuran | 83704-31-8 |

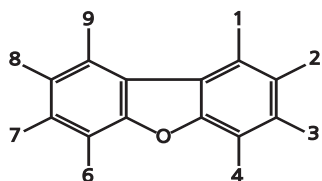
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZOFURANS

| ID Number* | Congener | CAS Number |
|------------|-----------------------------------|------------|
| 80 | 2,3,4,8-Tetrachlorodibenzofuran | 83704-32-9 |
| 81 | 2,3,6,7-Tetrachlorodibenzofuran | 57117-39-2 |
| 82 | 2,3,6,8-Tetrachlorodibenzofuran | 57117-37-0 |
| 83 | 2,3,7,8-Tetrachlorodibenzofuran | 51207-31-9 |
| 84 | 2,4,6,7-Tetrachlorodibenzofuran | 57117-38-1 |
| 85 | 2,4,6,8-Tetrachlorodibenzofuran | 58802-19-0 |
| 86 | 3,4,6,7-Tetrachlorodibenzofuran | 57117-40-6 |
| 87 | 1,2,3,4,6-Pentachlorodibenzofuran | 83704-47-6 |
| 88 | 1,2,3,4,7-Pentachlorodibenzofuran | 83704-48-7 |
| 89 | 1,2,3,4,8-Pentachlorodibenzofuran | 67517-48-0 |
| 90 | 1,2,3,4,9-Pentachlorodibenzofuran | 83704-49-8 |
| 91 | 1,2,3,6,7-Pentachlorodibenzofuran | 57117-42-7 |
| 92 | 1,2,3,6,8-Pentachlorodibenzofuran | 83704-51-2 |
| 93 | 1,2,3,6,9-Pentachlorodibenzofuran | 83704-52-3 |
| 94 | 1,2,3,7,8-Pentachlorodibenzofuran | 57117-41-6 |
| 95 | 1,2,3,7,9-Pentachlorodibenzofuran | 83704-53-4 |
| 96 | 1,2,3,8,9-Pentachlorodibenzofuran | 83704-54-5 |
| 97 | 1,2,4,6,7-Pentachlorodibenzofuran | 83704-50-1 |
| 98 | 1,2,4,6,8-Pentachlorodibenzofuran | 69698-57-3 |
| 99 | 1,2,4,6,9-Pentachlorodibenzofuran | 70648-24-7 |
| 100 | 1,2,4,7,8-Pentachlorodibenzofuran | 58802-15-6 |
| 101 | 1,2,4,7,9-Pentachlorodibenzofuran | 71998-74-8 |
| 102 | 1,2,4,8,9-Pentachlorodibenzofuran | 70648-23-6 |
| 103 | 1,2,6,7,8-Pentachlorodibenzofuran | 69433-00-7 |
| 104 | 1,2,6,7,9-Pentachlorodibenzofuran | 70872-82-1 |
| 105 | 1,3,4,6,7-Pentachlorodibenzofuran | 83704-36-3 |
| 106 | 1,3,4,6,8-Pentachlorodibenzofuran | 83704-55-6 |
| 107 | 1,3,4,6,9-Pentachlorodibenzofuran | 70648-15-6 |
| 108 | 1,3,4,7,8-Pentachlorodibenzofuran | 58802-16-7 |
| 109 | 1,3,4,7,9-Pentachlorodibenzofuran | 70648-20-3 |
| 110 | 1,3,6,7,8-Pentachlorodibenzofuran | 70648-21-4 |
| 111 | 1,4,6,7,8-Pentachlorodibenzofuran | 83704-35-2 |
| 112 | 2,3,4,6,7-Pentachlorodibenzofuran | 57117-43-8 |
| 113 | 2,3,4,6,8-Pentachlorodibenzofuran | 67481-22-5 |

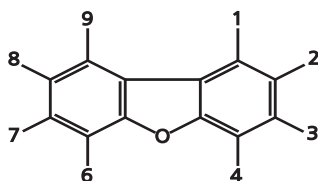
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED DIBENZOFURANS

| ID Number* | Congener | CAS Number |
|------------|---------------------------------------|------------|
| 114 | 2,3,4,7,8-Pentachlorodibenzofuran | 57117-31-4 |
| 115 | 1,2,3,4,6,7-Hexachlorodibenzofuran | 79060-60-9 |
| 116 | 1,2,3,4,6,8-Hexachlorodibenzofuran | 69698-60-8 |
| 117 | 1,2,3,4,6,9-Hexachlorodibenzofuran | 91538-83-9 |
| 118 | 1,2,3,4,7,8-Hexachlorodibenzofuran | 70648-26-9 |
| 119 | 1,2,3,4,7,9-Hexachlorodibenzofuran | 91538-84-0 |
| 120 | 1,2,3,4,8,9-Hexachlorodibenzofuran | 92341-07-6 |
| 121 | 1,2,3,6,7,8-Hexachlorodibenzofuran | 57117-44-9 |
| 122 | 1,2,3,6,7,9-Hexachlorodibenzofuran | 92341-06-5 |
| 123 | 1,2,3,6,8,9-Hexachlorodibenzofuran | |
| 124 | 1,2,3,7,8,9-Hexachlorodibenzofuran | 72918-21-9 |
| 125 | 1,2,4,6,7,8-Hexachlorodibenzofuran | 67562-40-7 |
| 126 | 1,2,4,6,7,9-Hexachlorodibenzofuran | 75627-02-0 |
| 127 | 1,2,4,6,8,9-Hexachlorodibenzofuran | 69698-59-5 |
| 128 | 1,3,4,6,7,8-Hexachlorodibenzofuran | 71998-75-9 |
| 129 | 1,3,4,6,7,9-Hexachlorodibenzofuran | 92341-05-4 |
| 130 | 2,3,4,6,7,8-Hexachlorodibenzofuran | 60851-34-5 |
| 131 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran | 67562-39-4 |
| 132 | 1,2,3,4,6,7,9-Heptachlorodibenzofuran | 70648-25-8 |
| 133 | 1,2,3,4,6,8,9-Heptachlorodibenzofuran | 69698-58-4 |
| 134 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran | 55673-89-7 |
| 135 | Octachlorodibenzofuran | 39001-02-0 |

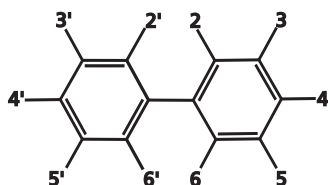
* Ballschmitter et. al.



SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|------------|--------------------------|------------|
| 1 | 2-Chlorobiphenyl | 2051-60-7 |
| 2 | 3-Chlorobiphenyl | 2051-61-8 |
| 3 | 4-Chlorobiphenyl | 2051-62-9 |
| 4 | 2,2'-Dichlorobiphenyl | 13029-08-8 |
| 5 | 2,3-Dichlorobiphenyl | 16605-91-7 |
| 6 | 2,3'-Dichlorobiphenyl | 25569-80-6 |
| 7 | 2,4-Dichlorobiphenyl | 33284-50-3 |
| 8 | 2,4'-Dichlorobiphenyl | 34883-43-7 |
| 9 | 2,5-Dichlorobiphenyl | 34883-39-1 |
| 10 | 2,6-Dichlorobiphenyl | 33146-45-1 |
| 11 | 3,3'-Dichlorobiphenyl | 2050-67-1 |
| 12 | 3,4-Dichlorobiphenyl | 2974-92-7 |
| 13 | 3,4'-Dichlorobiphenyl | 2974-90-5 |
| 14 | 3,5-Dichlorobiphenyl | 34883-41-5 |
| 15 | 4,4'-Dichlorobiphenyl | 2050-68-2 |
| 16 | 2,2',3-Trichlorobiphenyl | 38444-78-9 |
| 17 | 2,2',4-Trichlorobiphenyl | 37680-66-3 |
| 18 | 2,2',5-Trichlorobiphenyl | 37680-65-2 |
| 19 | 2,2',6-Trichlorobiphenyl | 38444-73-4 |
| 20 | 2,3,3'-Trichlorobiphenyl | 38444-84-7 |
| 21 | 2,3,4-Trichlorobiphenyl | 55702-46-0 |
| 22 | 2,3,4'-Trichlorobiphenyl | 38444-85-8 |
| 23 | 2,3,5-Trichlorobiphenyl | 55720-44-0 |
| 24 | 2,3,6-Trichlorobiphenyl | 55702-45-9 |
| 25 | 2,3',4-Trichlorobiphenyl | 55712-37-3 |
| 26 | 2,3',5-Trichlorobiphenyl | 38444-81-4 |
| 27 | 2,3',6-Trichlorobiphenyl | 38444-76-7 |
| 28 | 2,4,4'-Trichlorobiphenyl | 7012-37-5 |
| 29 | 2,4,5-Trichlorobiphenyl | 15862-07-4 |
| 30 | 2,4,6-Trichlorobiphenyl | 35693-92-6 |
| 31 | 2,4',5-Trichlorobiphenyl | 16606-02-3 |
| 32 | 2,4',6-Trichlorobiphenyl | 38444-77-8 |
| 33 | 2',3,4-Trichlorobiphenyl | 38444-86-9 |

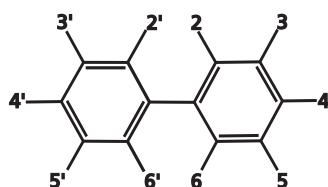
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|------------|-------------------------------|------------|
| 34 | 2',3,5-Trichlorobiphenyl | 37680-68-5 |
| 35 | 3,3',4-Trichlorobiphenyl | 37680-69-6 |
| 36 | 3,3',5-Trichlorobiphenyl | 38444-87-0 |
| 37 | 3,4,4'-Trichlorobiphenyl | 38444-90-5 |
| 38 | 3,4,5-Trichlorobiphenyl | 53555-66-1 |
| 39 | 3,4',5-Trichlorobiphenyl | 38444-88-1 |
| 40 | 2,2',3,3'-Tetrachlorobiphenyl | 38444-93-8 |
| 41 | 2,2',3,4-Tetrachlorobiphenyl | 52663-59-9 |
| 42 | 2,2',3,4'-Tetrachlorobiphenyl | 36559-22-5 |
| 43 | 2,2',3,5-Tetrachlorobiphenyl | 70362-46-8 |
| 44 | 2,2',3,5'-Tetrachlorobiphenyl | 41464-39-5 |
| 45 | 2,2',3,6-Tetrachlorobiphenyl | 70362-45-7 |
| 46 | 2,2',3,6'-Tetrachlorobiphenyl | 41464-47-5 |
| 47 | 2,2',4,4'-Tetrachlorobiphenyl | 2437-79-8 |
| 48 | 2,2',4,5-Tetrachlorobiphenyl | 70362-47-9 |
| 49 | 2,2',4,5'-Tetrachlorobiphenyl | 41464-40-8 |
| 50 | 2,2',4,6-Tetrachlorobiphenyl | 62796-65-0 |
| 51 | 2,2',4,6'-Tetrachlorobiphenyl | 68194-04-7 |
| 52 | 2,2',5,5'-Tetrachlorobiphenyl | 35693-99-3 |
| 53 | 2,2',5,6'-Tetrachlorobiphenyl | 41464-41-9 |
| 54 | 2,2',6,6'-Tetrachlorobiphenyl | 15968-05-5 |
| 55 | 2,3,3',4-Tetrachlorobiphenyl | 74338-24-2 |
| 56 | 2,3,3',4'-Tetrachlorobiphenyl | 41464-43-1 |
| 57 | 2,3,3',5-Tetrachlorobiphenyl | 70424-67-8 |
| 58 | 2,3,3',5'-Tetrachlorobiphenyl | 41464-49-7 |
| 59 | 2,3,3',6-Tetrachlorobiphenyl | 74472-33-6 |
| 60 | 2,3,4,4'-Tetrachlorobiphenyl | 33025-41-1 |
| 61 | 2,3,4,5-Tetrachlorobiphenyl | 33284-53-6 |
| 62 | 2,3,4,6-Tetrachlorobiphenyl | 54230-22-7 |
| 63 | 2,3,4',5-Tetrachlorobiphenyl | 74472-34-7 |
| 64 | 2,3,4',6-Tetrachlorobiphenyl | 52663-58-8 |
| 65 | 2,3,5,6-Tetrachlorobiphenyl | 33284-54-7 |
| 66 | 2,3',4,4'-Tetrachlorobiphenyl | 32598-10-0 |
| 67 | 2,3',4,5-Tetrachlorobiphenyl | 73575-53-8 |

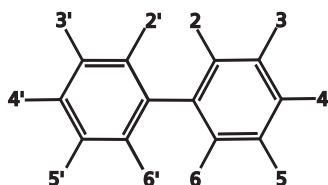
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|------------|---------------------------------|------------|
| 68 | 2,3',4,5'-Tetrachlorobiphenyl | 73575-52-7 |
| 69 | 2,3',4,6-Tetrachlorobiphenyl | 60233-24-1 |
| 70 | 2,3',4',5-Tetrachlorobiphenyl | 32598-11-1 |
| 71 | 2,3',4',6-Tetrachlorobiphenyl | 41464-46-4 |
| 72 | 2,3',5,5'-Tetrachlorobiphenyl | 41464-42-0 |
| 73 | 2,3',5',6-Tetrachlorobiphenyl | 74338-23-1 |
| 74 | 2,4,4',5-Tetrachlorobiphenyl | 32690-93-0 |
| 75 | 2,4,4',6-Tetrachlorobiphenyl | 32598-12-2 |
| 76 | 2',3,4,5-Tetrachlorobiphenyl | 70362-48-0 |
| 77 | 3,3',4,4'-Tetrachlorobiphenyl | 32598-13-3 |
| 78 | 3,3',4,5-Tetrachlorobiphenyl | 70362-49-1 |
| 79 | 3,3',4,5'-Tetrachlorobiphenyl | 41464-48-6 |
| 80 | 3,3',5,5'-Tetrachlorobiphenyl | 33284-52-5 |
| 81 | 3,4,4',5-Tetrachlorobiphenyl | 70362-50-4 |
| 82 | 2,2',3,3',4-Pentachlorobiphenyl | 52663-62-4 |
| 83 | 2,2',3,3',5-Pentachlorobiphenyl | 60145-20-2 |
| 84 | 2,2',3,3',6-Pentachlorobiphenyl | 52663-60-2 |
| 85 | 2,2',3,4,4'-Pentachlorobiphenyl | 65510-45-4 |
| 86 | 2,2',3,4,5-Pentachlorobiphenyl | 55312-69-1 |
| 87 | 2,2',3,4,5'-Pentachlorobiphenyl | 38380-02-8 |
| 88 | 2,2',3,4,6-Pentachlorobiphenyl | 55215-17-3 |
| 89 | 2,2',3,4,6'-Pentachlorobiphenyl | 73575-57-2 |
| 90 | 2,2',3,4',5-Pentachlorobiphenyl | 68194-07-0 |
| 91 | 2,2',3,4',6-Pentachlorobiphenyl | 68194-05-8 |
| 92 | 2,2',3,5,5'-Pentachlorobiphenyl | 52663-61-3 |
| 93 | 2,2',3,5,6-Pentachlorobiphenyl | 73575-56-1 |
| 94 | 2,2',3,5,6'-Pentachlorobiphenyl | 73575-55-0 |
| 95 | 2,2',3,5',6-Pentachlorobiphenyl | 38379-99-6 |
| 96 | 2,2',3,6,6'-Pentachlorobiphenyl | 73575-54-9 |
| 97 | 2,2',3',4,5-Pentachlorobiphenyl | 41464-51-1 |
| 98 | 2,2',3',4,6-Pentachlorobiphenyl | 60233-25-2 |
| 99 | 2,2',4,4',5-Pentachlorobiphenyl | 38380-01-7 |
| 100 | 2,2',4,4',6-Pentachlorobiphenyl | 39485-83-1 |
| 101 | 2,2',4,5,5'-Pentachlorobiphenyl | 37680-73-2 |

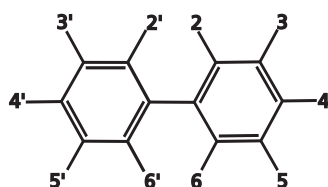
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|---------------|-----------------------------------|------------|
| 102 | 2,2',4,5,6'-Pentachlorobiphenyl | 68194-06-9 |
| 103 | 2,2',4,5',6-Pentachlorobiphenyl | 60145-21-3 |
| 104 | 2,2',4,6,6'-Pentachlorobiphenyl | 56558-16-8 |
| 105 | 2,3,3',4,4'-Pentachlorobiphenyl | 32598-14-4 |
| 106 | 2,3,3',4,5-Pentachlorobiphenyl | 70424-69-0 |
| 107 (BZ #108) | 2,3,3',4,5'-Pentachlorobiphenyl | 70362-41-3 |
| 108 (BZ #109) | 2,3,3',4,6-Pentachlorobiphenyl | 74472-35-8 |
| 109 (BZ #107) | 2,3,3',4',5-Pentachlorobiphenyl | 70424-68-9 |
| 110 | 2,3,3',4',6-Pentachlorobiphenyl | 38380-03-9 |
| 111 | 2,3,3',5,5'-Pentachlorobiphenyl | 39635-32-0 |
| 112 | 2,3,3',5,6-Pentachlorobiphenyl | 74472-36-9 |
| 113 | 2,3,3',5',6-Pentachlorobiphenyl | 68194-10-5 |
| 114 | 2,3,4,4',5-Pentachlorobiphenyl | 74472-37-0 |
| 115 | 2,3,4,4',6-Pentachlorobiphenyl | 74472-38-1 |
| 116 | 2,3,4,5,6-Pentachlorobiphenyl | 18259-05-7 |
| 117 | 2,3,4',5,6-Pentachlorobiphenyl | 68194-11-6 |
| 118 | 2,3',4,4',5-Pentachlorobiphenyl | 31508-00-6 |
| 119 | 2,3',4,4',6-Pentachlorobiphenyl | 56558-17-9 |
| 120 | 2,3',4,5,5'-Pentachlorobiphenyl | 68194-12-7 |
| 121 | 2,3',4,5',6-Pentachlorobiphenyl | 56558-18-0 |
| 122 | 2',3,3',4,5-Pentachlorobiphenyl | 76842-07-4 |
| 123 | 2',3,4,4',5-Pentachlorobiphenyl | 65510-44-3 |
| 124 | 2',3,4,5,5'-Pentachlorobiphenyl | 70424-70-3 |
| 125 | 2',3,4,5,6'-Pentachlorobiphenyl | 74472-39-2 |
| 126 | 3,3',4,4',5-Pentachlorobiphenyl | 57465-28-8 |
| 127 | 3,3',4,5,5'-Pentachlorobiphenyl | 39635-33-1 |
| 128 | 2,2',3,3',4,4'-Hexachlorobiphenyl | 38380-07-3 |
| 129 | 2,2',3,3',4,5-Hexachlorobiphenyl | 55215-18-4 |
| 130 | 2,2',3,3',4,5'-Hexachlorobiphenyl | 52663-66-8 |
| 131 | 2,2',3,3',4,6-Hexachlorobiphenyl | 61798-70-7 |
| 132 | 2,2',3,3',4,6'-Hexachlorobiphenyl | 38380-05-1 |
| 133 | 2,2',3,3',5,5'-Hexachlorobiphenyl | 35694-04-3 |
| 134 | 2,2',3,3',5,6-Hexachlorobiphenyl | 52704-70-8 |
| 135 | 2,2',3,3',5,6'-Hexachlorobiphenyl | 52744-13-5 |

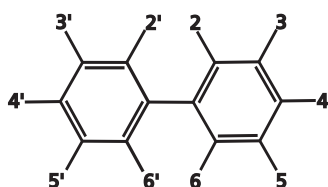
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|------------|-----------------------------------|------------|
| 136 | 2,2',3,3',6,6'-Hexachlorobiphenyl | 38411-22-2 |
| 137 | 2,2',3,4,4',5-Hexachlorobiphenyl | 35694-06-5 |
| 138 | 2,2',3,4,4',5'-Hexachlorobiphenyl | 35065-28-2 |
| 139 | 2,2',3,4,4',6-Hexachlorobiphenyl | 56030-56-9 |
| 140 | 2,2',3,4,4',6'-Hexachlorobiphenyl | 59291-64-4 |
| 141 | 2,2',3,4,5,5'-Hexachlorobiphenyl | 52712-04-6 |
| 142 | 2,2',3,4,5,6-Hexachlorobiphenyl | 41411-61-4 |
| 143 | 2,2',3,4,5,6'-Hexachlorobiphenyl | 68194-15-0 |
| 144 | 2,2',3,4,5',6-Hexachlorobiphenyl | 68194-14-9 |
| 145 | 2,2',3,4,6,6'-Hexachlorobiphenyl | 74472-40-5 |
| 146 | 2,2',3,4',5,5'-Hexachlorobiphenyl | 51908-16-8 |
| 147 | 2,2',3,4',5,6-Hexachlorobiphenyl | 68194-13-8 |
| 148 | 2,2',3,4',5,6'-Hexachlorobiphenyl | 74472-41-6 |
| 149 | 2,2',3,4',5',6-Hexachlorobiphenyl | 38380-04-0 |
| 150 | 2,2',3,4',6,6'-Hexachlorobiphenyl | 68194-08-1 |
| 151 | 2,2',3,5,5',6-Hexachlorobiphenyl | 52663-63-5 |
| 152 | 2,2',3,5,6,6'-Hexachlorobiphenyl | 68194-09-2 |
| 153 | 2,2',4,4',5,5'-Hexachlorobiphenyl | 35065-27-1 |
| 154 | 2,2',4,4',5,6'-Hexachlorobiphenyl | 60145-22-4 |
| 155 | 2,2',4,4',6,6'-Hexachlorobiphenyl | 33979-03-2 |
| 156 | 2,3,3',4,4',5-Hexachlorobiphenyl | 38380-08-4 |
| 157 | 2,3,3',4,4',5'-Hexachlorobiphenyl | 69782-90-7 |
| 158 | 2,3,3',4,4',6-Hexachlorobiphenyl | 74472-42-7 |
| 159 | 2,3,3',4,5,5'-Hexachlorobiphenyl | 39635-35-3 |
| 160 | 2,3,3',4,5,6-Hexachlorobiphenyl | 41411-62-5 |
| 161 | 2,3,3',4,5',6-Hexachlorobiphenyl | 74472-43-8 |
| 162 | 2,3,3',4',5,5'-Hexachlorobiphenyl | 39635-34-2 |
| 163 | 2,3,3',4',5,6-Hexachlorobiphenyl | 74472-44-9 |
| 164 | 2,3,3',4',5',6-Hexachlorobiphenyl | 74472-45-0 |
| 165 | 2,3,3',5,5',6-Hexachlorobiphenyl | 74472-46-1 |
| 166 | 2,3,4,4',5,6-Hexachlorobiphenyl | 41411-63-6 |
| 167 | 2,3',4,4',5,5'-Hexachlorobiphenyl | 52663-72-6 |
| 168 | 2,3',4,4',5',6-Hexachlorobiphenyl | 59291-65-5 |
| 169 | 3,3',4,4',5,5'-Hexachlorobiphenyl | 32774-16-6 |

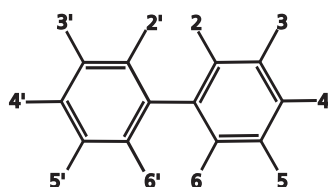
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|---------------|--|------------|
| 170 | 2,2',3,3',4,4',5-Heptachlorobiphenyl | 35065-30-6 |
| 171 | 2,2',3,3',4,4',6-Heptachlorobiphenyl | 52663-71-5 |
| 172 | 2,2',3,3',4,5,5'-Heptachlorobiphenyl | 52663-74-8 |
| 173 | 2,2',3,3',4,5,6-Heptachlorobiphenyl | 68194-16-1 |
| 174 | 2,2',3,3',4,5,6'-Heptachlorobiphenyl | 38411-25-5 |
| 175 | 2,2',3,3',4,5',6-Heptachlorobiphenyl | 40186-70-7 |
| 176 | 2,2',3,3',4,6,6'-Heptachlorobiphenyl | 52663-65-7 |
| 177 | 2,2',3,3',4',5,6-Heptachlorobiphenyl | 52663-70-4 |
| 178 | 2,2',3,3',5,5',6-Heptachlorobiphenyl | 52663-67-9 |
| 179 | 2,2',3,3',5,6,6'-Heptachlorobiphenyl | 52663-64-6 |
| 180 | 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 35065-29-3 |
| 181 | 2,2',3,4,4',5,6-Heptachlorobiphenyl | 74472-47-2 |
| 182 | 2,2',3,4,4',5,6'-Heptachlorobiphenyl | 60145-23-5 |
| 183 | 2,2',3,4,4',5',6-Heptachlorobiphenyl | 52663-69-1 |
| 184 | 2,2',3,4,4',6,6'-Heptachlorobiphenyl | 74472-48-3 |
| 185 | 2,2',3,4,5,5',6-Heptachlorobiphenyl | 52712-05-7 |
| 186 | 2,2',3,4,5,6,6'-Heptachlorobiphenyl | 74472-49-4 |
| 187 | 2,2',3,4',5,5',6-Heptachlorobiphenyl | 52663-68-0 |
| 188 | 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 74487-85-7 |
| 189 | 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 39635-31-9 |
| 190 | 2,3,3',4,4',5,6-Heptachlorobiphenyl | 41411-64-7 |
| 191 | 2,3,3',4,4',5',6-Heptachlorobiphenyl | 74472-50-7 |
| 192 | 2,3,3',4,5,5',6-Heptachlorobiphenyl | 74472-51-8 |
| 193 | 2,3,3',4',5,5',6-Heptachlorobiphenyl | 69782-91-8 |
| 194 | 2,2',3,3',4,4',5,5'-Octachlorobiphenyl | 35694-08-7 |
| 195 | 2,2',3,3',4,4',5,6-Octachlorobiphenyl | 52663-78-2 |
| 196 | 2,2',3,3',4,4',5,6'-Octachlorobiphenyl | 42740-50-1 |
| 197 | 2,2',3,3',4,4',6,6'-Octachlorobiphenyl | 33091-17-7 |
| 198 | 2,2',3,3',4,5,5',6-Octachlorobiphenyl | 68194-17-2 |
| 199 (BZ #201) | 2,2',3,3',4,5,5',6'-Octachlorobiphenyl | 52663-75-9 |
| 200 (BZ #199) | 2,2',3,3',4,5,6,6'-Octachlorobiphenyl | 52663-73-7 |
| 201 (BZ #200) | 2,2',3,3',4,5',6,6'-Octachlorobiphenyl | 40186-71-8 |
| 202 | 2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 2136-99-4 |
| 203 | 2,2',3,4,4',5,5',6-Octachlorobiphenyl | 52663-76-0 |

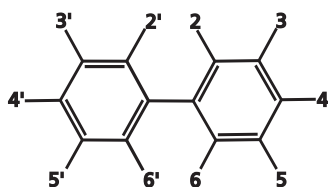
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SYSTEMATIC NUMBERING OF CHLORINATED BIPHENYLS

| ID Number* | Congener | CAS Number |
|------------|--|------------|
| 204 | 2,2',3,4,4',5,6,6'-Octachlorobiphenyl | 74472-52-9 |
| 205 | 2,3,3',4,4',5,5',6-Octachlorobiphenyl | 74472-53-0 |
| 206 | 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 40186-72-9 |
| 207 | 2,2',3,3',4,4',5,6,6'-Nonachlorobiphenyl | 52663-79-3 |
| 208 | 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 52663-77-1 |
| 209 | Decachlorobiphenyl | 2051-24-3 |

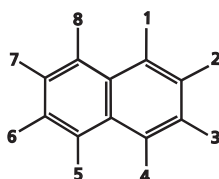
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SYSTEMATIC NUMBERING OF CHLORINATED NAPHTHALENES

| ID Number* | Congener | CAS Number |
|------------|----------------------------|------------|
| 1 | 1-Chloronaphthalene | 90-13-1 |
| 2 | 2-Chloronaphthalene | 91-58-7 |
| 3 | 1,2-Dichloronaphthalene | 2050-69-3 |
| 4 | 1,3-Dichloronaphthalene | 2198-75-6 |
| 5 | 1,4-Dichloronaphthalene | 1825-31-6 |
| 6 | 1,5-Dichloronaphthalene | 1825-30-5 |
| 7 | 1,6-Dichloronaphthalene | 2050-72-8 |
| 8 | 1,7-Dichloronaphthalene | 2050-73-9 |
| 9 | 1,8-Dichloronaphthalene | 2050-74-0 |
| 10 | 2,3-Dichloronaphthalene | 2050-75-1 |
| 11 | 2,6-Dichloronaphthalene | 2065-70-5 |
| 12 | 2,7-Dichloronaphthalene | 2198-77-8 |
| 13 | 1,2,3-Trichloronaphthalene | 50402-52-3 |
| 14 | 1,2,4-Trichloronaphthalene | 50402-51-2 |

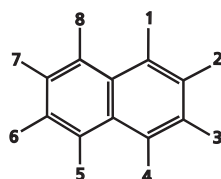
* Wiedmann and Ballschmitter



SYSTEMATIC NUMBERING OF CHLORINATED NAPHTHALENES

| ID Number* | Congener | CAS Number |
|------------|--------------------------------|-------------|
| 15 | 1,2,5-Trichloronaphthalene | 55720-33-7 |
| 16 | 1,2,6-Trichloronaphthalene | 51570-44-6 |
| 17 | 1,2,7-Trichloronaphthalene | 55720-34-8 |
| 18 | 1,2,8-Trichloronaphthalene | 55720-35-9 |
| 19 | 1,3,5-Trichloronaphthalene | 51570-43-5 |
| 20 | 1,3,6-Trichloronaphthalene | 55720-36-0 |
| 21 | 1,3,7-Trichloronaphthalene | 55720-37-1 |
| 22 | 1,3,8-Trichloronaphthalene | 55720-38-2 |
| 23 | 1,4,5-Trichloronaphthalene | 2437-55-0 |
| 24 | 1,4,6-Trichloronaphthalene | 2737-54-9 |
| 25 | 1,6,7-Trichloronaphthalene | 55720-39-3 |
| 26 | 2,3,6-Trichloronaphthalene | 55720-40-6 |
| 27 | 1,2,3,4-Tetrachloronaphthalene | 20020-02-4 |
| 28 | 1,2,3,5-Tetrachloronaphthalene | 53555-63-8 |
| 29 | 1,2,3,6-Tetrachloronaphthalene | |
| 30 | 1,2,3,7-Tetrachloronaphthalene | 55720-41-7 |
| 31 | 1,2,3,8-Tetrachloronaphthalene | 149864-81-3 |
| 32 | 1,2,4,5-Tetrachloronaphthalene | 6733-54-6 |
| 33 | 1,2,4,6-Tetrachloronaphthalene | 51570-45-7 |
| 34 | 1,2,4,7-Tetrachloronaphthalene | 67922-21-8 |
| 35 | 1,2,4,8-Tetrachloronaphthalene | 6529-87-9 |
| 36 | 1,2,5,6-Tetrachloronaphthalene | 67922-22-9 |
| 37 | 1,2,5,7-Tetrachloronaphthalene | 67922-23-0 |
| 38 | 1,2,5,8-Tetrachloronaphthalene | 149864-80-2 |
| 39 | 1,2,6,7-Tetrachloronaphthalene | 149864-79-9 |
| 40 | 1,2,6,8-Tetrachloronaphthalene | 67922-24-1 |
| 41 | 1,2,7,8-Tetrachloronaphthalene | 149864-82-4 |
| 42 | 1,3,5,7-Tetrachloronaphthalene | 53555-64-9 |
| 43 | 1,3,5,8-Tetrachloronaphthalene | 31604-28-1 |
| 44 | 1,3,6,7-Tetrachloronaphthalene | 55720-42-8 |
| 45 | 1,3,6,8-Tetrachloronaphthalene | 150224-15-0 |
| 46 | 1,4,5,8-Tetrachloronaphthalene | 3432-57-3 |
| 47 | 1,4,6,7-Tetrachloronaphthalene | 55720-43-9 |
| 48 | 2,3,6,7-Tetrachloronaphthalene | |

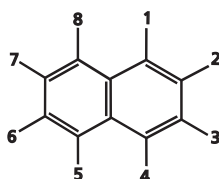
* Wiedmann and Ballschmitter



SYSTEMATIC NUMBERING OF CHLORINATED NAPHTHALENES

| ID Number* | Congener | CAS Number |
|------------|--------------------------------------|-------------|
| 49 | 1,2,3,4,5-Pentachloronaphthalene | 67922-25-2 |
| 50 | 1,2,3,4,6-Pentachloronaphthalene | 67922-26-3 |
| 51 | 1,2,3,5,6-Pentachloronaphthalene | |
| 52 | 1,2,3,5,7-Pentachloronaphthalene | 53555-65-0 |
| 53 | 1,2,3,5,8-Pentachloronaphthalene | 150224-24-1 |
| 54 | 1,2,3,6,7-Pentachloronaphthalene | 150224-16-1 |
| 55 | 1,2,3,6,8-Pentachloronaphthalene | 150224-23-0 |
| 56 | 1,2,3,7,8-Pentachloronaphthalene | 150205-21-3 |
| 57 | 1,2,4,5,6-Pentachloronaphthalene | 150224-20-7 |
| 58 | 1,2,4,5,7-Pentachloronaphthalene | 150224-19-4 |
| 59 | 1,2,4,5,8-Pentachloronaphthalene | 150224-25-2 |
| 60 | 1,2,4,6,7-Pentachloronaphthalene | 150224-17-2 |
| 61 | 1,2,4,6,8-Pentachloronaphthalene | 150224-22-9 |
| 62 | 1,2,4,7,8-Pentachloronaphthalene | |
| 63 | 1,2,3,4,5,6-Hexachloronaphthalene | 58877-88-6 |
| 64 | 1,2,3,4,5,7-Hexachloronaphthalene | 67927-27-4 |
| 65 | 1,2,3,4,5,8-Hexachloronaphthalene | 103426-93-3 |
| 66 | 1,2,3,4,6,7-Hexachloronaphthalene | 103426-96-6 |
| 67 | 1,2,3,5,6,7-Hexachloronaphthalene | 103426-97-7 |
| 68 | 1,2,3,5,6,8-Hexachloronaphthalene | 103426-95-5 |
| 69 | 1,2,3,5,7,8-Hexachloronaphthalene | 103426-94-4 |
| 70 | 1,2,3,6,7,8-Hexachloronaphthalene | |
| 71 | 1,2,4,5,6,8-Hexachloronaphthalene | 90948-28-0 |
| 72 | 1,2,4,5,7,8-Hexachloronaphthalene | 103426-92-2 |
| 73 | 1,2,3,4,5,6,7-Heptachloronaphthalene | 58863-14-2 |
| 74 | 1,2,3,4,5,6,8-Heptachloronaphthalene | 58863-15-3 |
| 75 | Octachloronaphthalene | 2234-13-1 |

* Wiedmann and Ballschmitter





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AR-1523

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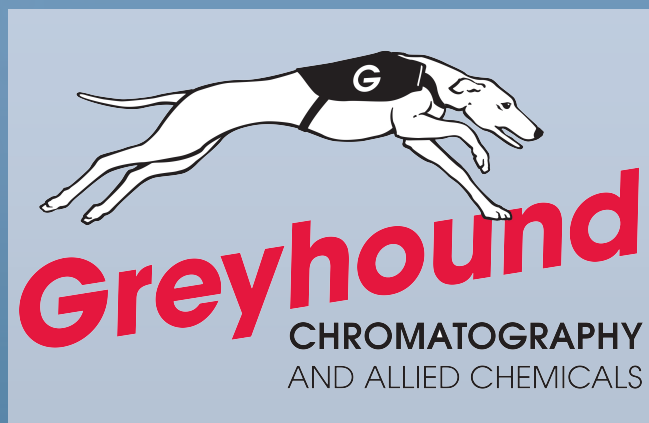
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