



Cambridge Isotope
Laboratories, Inc.

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Standards for Environmental, Food, Water and Exposure Analysis



Solutions for a Greener World



Environmental Contaminant Standards from CIL

Joel Bradley, PhD
Chief Executive Officer

Cambridge Isotope Laboratories, Inc. (CIL) is pleased to present its newest catalog of isotopically labeled and unlabeled standards designed to optimize environmental, food, water, and exposure analyses.

For the past 35 years, CIL has been the world's leading supplier of isotope-labeled standards to leading environmental analytical laboratories. In that time there have been substantial changes in the nature of analytical standards, and the nature of stakeholders who rely on isotopically labeled standards for rigorous analysis using isotope dilution mass spectrometry (IDMS).

Perhaps the most significant change has been in the markets CIL serves. CIL's customer base was once dominated by a handful of labs using high-resolution gas chromatography–mass spectrometry (HRGC-MS) to measure, with increasingly lower quantitation limits, the presence of dioxins, PCBs, priority pollutants, and legacy pesticides such as DDT in soil, water, air, and biota. These important analyses continue, but the scope of work has expanded rapidly into areas such as food, drinking water, and human-exposure testing. Target analytes have expanded in all directions to include compounds that were of interest to only a few visionaries in the early 1980s. Flame retardants, organophosphate (OP) pesticides, and pharmaceutical and personal care products (PPCPs), just to name a few, were just coming on to the radar at the turn of the millennium and now represent a significant portion of CIL's product offerings in the trace and ultratrace analytical arena.

CIL's latest efforts include many novel developments on compounds that have been targeted for decades, such as polycyclic aromatic compounds (PACs) and polychloronaphthalenes (PCNs), as well as compounds of emerging concern such as perfluorinated compounds (PFCs) and bisphenols (A, S, F, AF, E, P, AP and Z). CIL has also been developing many isotopically labeled standards for new-use pesticides, including several key neonicotinoid insecticides that are under investigation for potential contribution to colony collapse disorder (CCD) in honey bee populations, fipronil and pyrethroids used in indoor pest-control applications, and metabolites, byproducts, and decomposition products that are potential biomarkers.

So whether you analyze legacy POPs compounds or are studying compounds of emerging concern, CIL is committed to producing the critical standards needed by the global analytical community to discover, track, quantify, and, ultimately, help resolve contamination and exposure problems.

Joel C. Bradley, PhD
Chief Executive Officer

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A Few Words from the Environmental Products Team

Cambridge Isotope Laboratories, Inc. (CIL) is very pleased to present its latest catalog of "Standards for Environmental, Food, Water, and Exposure Analysis." As you may have noticed, the first new change in this catalog is the title itself. CIL recognizes that many organic contaminants are not only environmental pollutants, but may also be food and water contaminants, and ultimately compounds of concern in human exposure studies. As our product lines continue to grow, we find our standards being used not only as "environmental" contaminant standards, but increasingly in these interrelated applications and sample matrices. So whether you are testing PCBs in sediment, pesticides or dioxins and furans in food and feed, PPCPs in industrial effluent, flame retardants in house dust, or PAHs and tobacco metabolites in biomonitoring studies, you will find an extensive set of standards in the following pages to meet your analytical needs.

As new technologies and applications to study environmental, food, water, and exposure contaminants advance, CIL continues to maintain a leadership role working with researchers to develop new standards to meet the latest needs. With new instrumentation and extraction procedures allowing detection limits approaching attogram levels in some cases, the need for highly accurate isotope standards has never been greater. In this catalog you will find approximately 600 new standards for legacy and emerging compounds. A partial list of the recently developed projects includes:

- New combined dioxin/furan/PCB mixtures
- New ¹³C-labeled PAHs, PAH metabolites, and PAH standard mixtures
- Expanded offering of ¹³C-labeled and unlabeled PCNs and PCN mixtures
- Many new ¹³C-labeled pesticides, including fipronil and neonicotinoid insecticides
- Augmented list of toxaphene congeners, and new EPA Method 8276 standard mixtures
- Expanded set of tobacco-specific nitrosamines
- PFOS and other PFCs
- New ¹³C-labeled and unlabeled cyclic methyl siloxanes
- Bisphenol S and other bisphenol analogs
- Expanded phthalate monoester offerings
- New brominated, chlorinated, and phosphate flame retardants
- ¹³C-labeled short-chain chlorinated paraffins
- Various biomarker compounds and metabolites for exposure analysis

You can find all of this valuable information and more online as well. CIL launched a new website in 2014 to provide a more user-friendly experience when searching for product information, placing orders and quote requests, and keeping up to date with the latest news and information from CIL. As part of the interactive approach, visitors to our website can immediately access updated product information and documentation, such as a certificate of analysis (CoA) and safety data sheets (SDS), and can also find real-time product availability as part of our e-commerce functionality. An online version of this catalog can be accessed in the "Literature Library" section of the "Sales and Technical Support" tab. The Literature Library also includes newsletters, application notes, and new product "Spotlights." Visit www.isotope.com to see more.

So please take a look through our new catalog and enjoy the information provided, as well as our newly expanded listings of "Standards for Environmental, Food, Water, and Exposure Analysis!"

Environmental Products Team



Ben Priest
Business Development
Manager



Terry Grim
Regional Sales
Manager – North
America, Latin America
and Japan



Kayla Meehan
Sales and Marketing
Coordinator

Corporate Overview

CIL is the world leader in the separation and manufacture of stable (nonradioactive) isotopes and isotope-labeled compounds.

With over 400 employees and laboratories in four countries, CIL specializes in the process of labeling biochemical and organic compounds with highly enriched stable isotopes of carbon, hydrogen, nitrogen and oxygen. Our chemists substitute a common atom for a rare, highly valued isotopic component so that the final product can be readily measured or traced using mass spectrometry (MS) or nuclear magnetic resonance (NMR). CIL's products are utilized in laboratories, health care facilities, and medical, government and academic research centers worldwide. We are proud that CIL products have contributed to medical advancements in cancer research, new drug development, environmental and food analysis, genomics and proteomics, and medical diagnostic research.

CIL's vision began when it was founded in 1981 by Dr. Joel Bradley, an organic chemist from MIT. Drawing on a commitment to high-quality products, superior customer service, innovative new products and breadth of product lines, CIL quickly emerged as a leader in its field. CIL now produces more than 15,000 products and has ISO 13485, ISO/IEC 17025 and ISO Guide 34 quality systems, as well as cGMP production capabilities. The CIL group is comprised of five companies: Cambridge Isotope Laboratories, Inc. (CIL) and CIL Isotope Separations (CIS) in the United States; CIL Canada, Inc. in Montreal, Canada; Euriso-Top in Saclay, France; and ABX GmbH in Dresden, Germany.

CIL has worked closely with industry leaders and researchers to provide stable isotope-labeled tools needed for improved quantitation of complex systems. This has been particularly true in the last decade, when many innovative techniques for determining biomarkers for the presence, progression and monitoring of therapeutic response have emerged from the fields of MS-based proteomics and metabolomics.

CIL takes great pride in being able to offer a wide range of NMR solvents with the highest isotopic enrichment and chemical purity. All NMR solvents undergo thorough quality-control testing during the manufacturing and packaging process to verify that product quality is maintained. Continuous improvement and high-quality standards make CIL the preferred supplier to synthesis groups and CROs/CMOs worldwide.

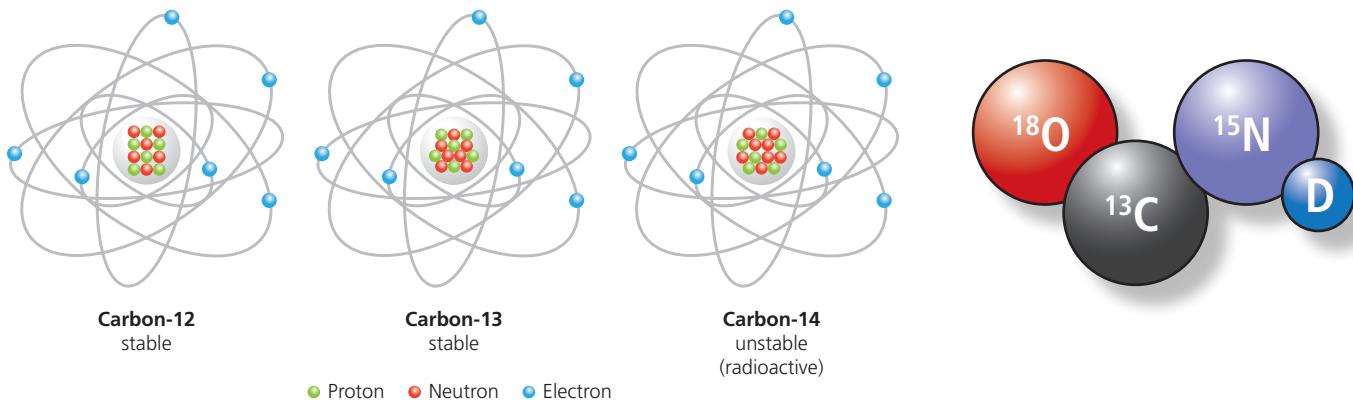
For more than 35 years, CIL has provided stable isotope-labeled standards for the analysis of organic contaminants using isotope dilution mass spectrometry (IDMS). The use of IDMS is widely accepted as an extremely accurate and robust analytical technique, yielding data of the highest quality. Through the years, CIL collaborations with leading researchers have led to development of novel products to help address new environment issues. From the early 1980s working with government laboratories to develop the first commercially available standards for use in the accurate determination of dioxins and furans in people exposed to Agent Orange, to the mid-1990s collaborating with groups around the world to develop new brominated diphenyl ether flame-retardant standards, to the recent development of labeled standards for fipronil and neonicotinoid insecticides to assist researchers studying bee colony collapse disorder, CIL continually assists researchers studying important global issues.

Dr. Bradley and the CIL executive team all share the same commitment to quality and service. CIL's experts collaborate with all of their customers to aid in pivotal research that is being conducted in laboratories worldwide. Our partnerships not only help to support our global reach, but allow us to bring forward innovative products to aid our customers' pursuit of scientific discovery.

What Is an Isotope?

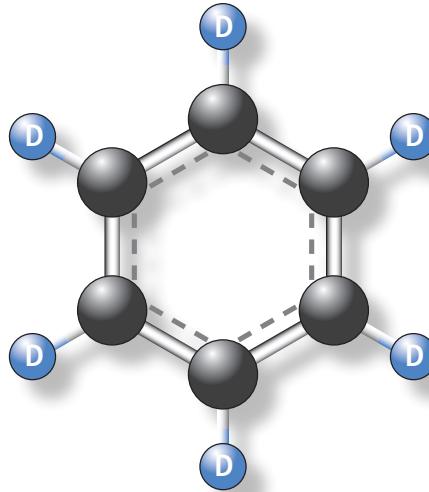
An isotope is any of two or more forms of a chemical element that has a different number of neutrons in the nucleus. There are 275 stable isotopes of the 81 nonradioactive elements, in

addition to over 800 radioactive isotopes. Every element has known isotopic forms. Isotopes of a single element possess almost identical chemical properties.



Isotopic Enrichment

Isotopic enrichment is the average enrichment for each labeled atom in the molecule. It is not the percentage of molecules that are completely isotopically labeled. For instance, benzene (D_6 , 99%) is not 99% C_6D_6 and 1% C_6H_6 . Each of the six hydrogen atoms has a 99% chance of being a deuterium atom ($^2H = D$), and a 1% chance of being protium (1H). Thus, $(99\%)^6$, or about 94% of the benzene molecules, will have a molecular mass that is six atomic mass units (amu) higher than native (unlabeled) benzene. About 6% will have a molecular mass that is 5 amu higher than native benzene. Theoretically, only $(1\%)^6$, or about $10^{-10}\%$, will have the molecular mass of native benzene.



Cambridge Isotope Laboratories, Inc. Facilities

CIL has state-of-the-art production facilities for cGMP and non-cGMP manufacturing at its locations in Andover and Tewksbury, Massachusetts.

CIL World Headquarters and cGMP Production Laboratories Tewksbury, MA USA

CIL moved into its new Tewksbury, Massachusetts, facility in the spring of 2013. As the new corporate headquarters, this facility houses the executive team, as well as sales, marketing, finance, regulatory affairs and cGMP production staff. In addition to corporate office space, the facility has a state-of-the-art cGMP suite, which includes production laboratories, dedicated isolation rooms, a dedicated analytical laboratory, a packaging laboratory and a development laboratory.



CIL Production Laboratories Andover, MA USA

CIL's primary production facility in Andover, Massachusetts, is dedicated to the manufacture of deuterated NMR solvents, stable isotope-labeled chemicals and gases, as well as specific cGMP products. This facility is home to operations staff and production and quality-control teams.

The formulations group has over 30 years' experience formulating highly purified labeled materials into high-quality quantitative solutions as analytical standards, either as single-component products or multi-component mixes and calibration solutions.

The quality-control lab is equipped with a wide array of instrumentation, including gas chromatograph/mass spectrometers (GC/MS), high-field NMRs, HPLCs and an FT-IR. CIL's chemistry laboratories are equipped with apparatus for both large-scale (50+ liters) and microscale chemistry, which includes equipment for high-pressure gas reactions, pH and temperature-controlled enzyme chemistry, high-resolution distillation processes, and catalytic reduction with both hydrogen and deuterium. The production laboratories are also equipped with analytical equipment for in-process testing, including GC-FID, GC-ECD and HPLC with UV, RI, ELSD and MS detectors. All of these resources allow CIL to consistently produce products with high chemical and isotopic purity.





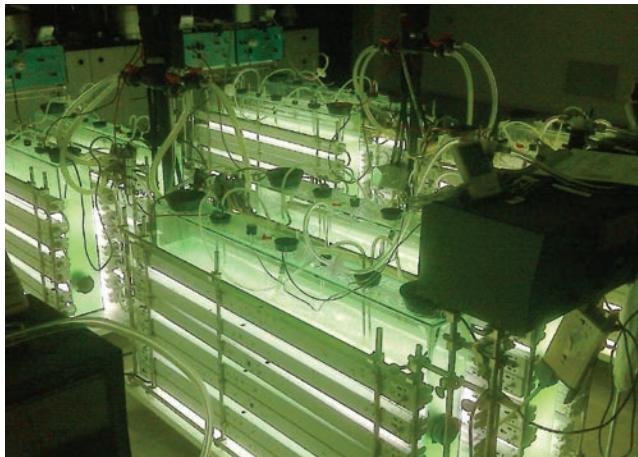
CIL Isotope Separations, LLC (CIS) Xenia, OH USA

CIL is the world leader in the separation of ^{13}C and ^{18}O . CIL separates both ^{13}C and ^{18}O at its Xenia, OH, facility and has the world's largest production capacity for both ^{13}C and ^{18}O . CIL also has the only nongovernmental D₂O enrichment columns in the world located at its CIS facility.



Euriso-Top Saclay, France

Euriso-Top was founded in January 1991 by a group of researchers from the Commissariat à l'Energie Atomique (CEA). Euriso-Top is Europe's leading producer of deuterated NMR solvents, cGMP urea and stable isotope-labeled compounds. Its quality control and production laboratories are equipped with NMR, MS, HPLC, GC, IR and UV instruments.



CIL Canada, Inc. Montreal, Canada

CIL Canada, Inc. is CIL's biotech laboratory facility, which produces carbohydrates, enriched media and amino acids for drug-discovery applications. CIL Canada specializes in algal biosynthesis, including spirulina, chlorella and a variety of other algal strains for NMR and proteomics applications.



ABX GmbH Dresden, Germany

ABX is the world's leading supplier of ^{18}F positron emission tomography (PET) precursors, reagent kits and cassettes, including, but not limited to, kits for FDG, FLT, F-choline, NaF, F-Miso and FET. ABX specializes in the manufacture and development of chemicals for nuclear medicine, and its cGMP-approved laboratories, class 100 clean rooms and cGMP radiochemistry development hot lab uniquely position ABX to provide complete PET and SPECT chemistry solutions to radiochemists and radiopharmacists worldwide. ABX's radiochemistry hot lab is equipped with most of the leading commercial PET tracer synthesis boxes and allows ABX to assist customers with the optimization and development of new tracers.

Ordering Information

Placing an Order

Phone: 1.800.322.1174 (North America) or
1.978.749.8000 (International)
Office hours are 8:00 a.m. to 5:30 p.m.
Eastern Standard Time (EST)

Fax: 1.978.749.2768

Email: envsales@isotope.com (North America)
intlsales@isotope.com (International)

E-commerce: Visit <http://shop.isotope.com> to request
a quote, place orders, obtain product
information or submit technical questions.

CIL products are continually updated on the website so be
sure to visit <http://shop.isotope.com> for current information.

Please help us to expedite the shipment of your order by
including the following information:

- Shipping address, including street
- Billing address
- Purchase order number or credit card information
- CIL catalog number and product name
- Quantity: mg (milligrams), g (grams), kg (kilograms),
mL (milliliters), L (liters), etc., as applicable, including
number of units
- Catalog price or CIL quotation number with date given
- Special instructions for packaging or shipping
- Your name, phone number and email address
- End user name, phone number and email address
(if different)
- Preferred mode of shipping (e.g. FedEx or UPS)
- \$50 minimum order

We do not require written confirmation of phone orders
for established customers.

First-Time Orders

If ordering for the first time, please email or fax the following
information on company letterhead to establish a line of
credit with a copy of your order:

- A federal tax identification number
- Three credit/banking references

Also include your shipping address, billing address, phone,
fax, email and URL address.

*To expedite delivery of your first order, prepayment should
be made by credit card or wire transfer in US funds.*

Pricing Information and Terms of Sale

North American Orders

- All prices are in US dollars. Any importation costs for
international orders are not included. Please consult our
Customer Service Department for pricing information
or packaging options.
- When stock is available and subdivision is possible, we will
accept orders for smaller than catalog amounts. Please
request a quotation as a quantity discount may apply.
- Please note that prices are subject to change without
notice. Occasionally the inventory of some products
listed may become depleted. Replacement of stock
may be subject to a minimum order quantity.
- You may check stock and confirm prices by contacting
the CIL Customer Service Department at 1.800.322.1174
(North America only) or envsales@isotope.com.
- CIL will be pleased to assist customers with firm written
quotations. Most quotes are valid for 30-60 days. Longer
terms may be granted by CIL upon request.
- Net 30 days from invoice date with prior credit approval.
Past-due invoices will be subject to a 1.5% per month
service charge; 18% per annum. We reserve the right
to request payment in advance or COD terms on initial
orders with CIL.
- We also accept VISA, MasterCard, American Express and
university purchasing card orders.
- Shipping terms are FCA Andover, MA USA. Any damage
to the package or product in transit is the buyer's
responsibility to adjust with the carrier.
- Domestic shipping charges will be added to invoices
(unless collect shipment is requested).

International Orders

- CIL has an extensive international sales network of
over 33 representatives in 27 different countries.
- For international orders or quotations, please contact
CIL International Sales at intlsales@isotope.com or
+1.978.749.8000. For a complete distributor listing,
please visit isotope.com.
- Representatives and agents are available to assist you
with your requirements for our products. Please consult
your local CIL representative for appropriate pricing and
payment terms. Shipping charges and any applicable
import duties and taxes will be added to orders placed
with distributors.
- For direct orders, CIL generally requires prepayment in
US dollars by an international bank check or bank wire
transfer. We will be pleased to provide *pro forma* invoices
upon request. Shipping charges will be added to direct
orders. Any applicable import duties and taxes will be
charged to the purchaser by the shipping company or
customs agent.
- Shipping terms are FCA Andover, MA USA. Any damage to
the package or product in transit is the buyer's responsibility
to adjust with the carrier.

Shipping Information

USA

- Shipments within the United States will be sent via UPS, FedEx, or truck.
- Orders within the United States for in-stock items placed before 2 p.m. EST can ship the same day via FedEx.

Canada

- Canadian shipments will be sent via FedEx or truck.
- Please include the name of your customs broker.
- Orders to Canada for in-stock items will ship one to two working days after receipt of purchase order.
- There are handling charges for orders of hazardous materials shipped via FedEx.

International

- International shipments will be sent via FedEx or best method.
- CIL tries to be as cost effective as possible, but the carrier may assess additional charges.

We will accommodate your shipping instructions whenever it is feasible to do so. CIL reserves the right to change the method of transportation, if required, to comply with transportation regulations. Such a change would not alter your responsibility for payment of shipping charges.

Additional shipping charges may apply.

Return Shipment Policy

Returns may be made within 30 days of shipment with prior approval from CIL. We reserve the right to impose restocking charges when a return is at the sole option of the buyer.

The buyer is responsible for approving the quality and quantity of any product within the 30-day period stated above. If an error by CIL results in an incorrect or duplicate shipment, a replacement will be sent or the appropriate credit allowed. We typically request return of the original product. Product returns must reference the original purchase order number, CIL order number (e.g. DB-A1000), Returned Goods Authorization (RGA) number, and the date CIL authorized the return. Under no circumstances will credit or replacement be given for products without prior authorization by CIL.

Product Information

Documentation

A Certificate of Analysis (COA) and a Safety Data Sheet (SDS) are supplied with every shipment. Additional product information may be available upon request.

The chemical purity (CP) specification for CIL products is 98% unless otherwise indicated.

Limited Warranty

CIL represents that the products are, as of the date of shipment, as described in CIL's applicable product literature. CIL makes no other warranty, express or implied, with respect to its products, including any warranty of merchantability or fitness for any particular purpose. CIL's maximum liability for any reason shall be to replace any nonconforming product or refund the applicable purchase price.

Environmental Food, Water, and Exposure Products Information

There are potential hazards associated with the use of any chemical, and CIL's environmental, food, water and exposure analysis standards may require additional considerations. Customers are encouraged to consult standard safety references for the proper use and handling of CIL products. While every effort was made to ensure the information in this catalog is correct, users of CIL products are responsible for confirming product information.

Neat Standards

Neat/crystalline standards contain approximately the stated mass, as crystal sizes and static electricity may make it difficult to weigh exact milligram quantities. Stated weights should not be used to prepare quantitative standards. Material should be weighed prior to standard formulation. Small amounts may need to be transferred with the use of solvent; the vial should be weighed before transfer, and after all solvent has been evaporated. During shipment, small but significant amounts of material may shift into the vial cap or ampoule tip. Traceable weighing records, or weights tailored to user-specified targets, can be obtained for a reasonable surcharge.

Quantitative Solutions

Quantitative standard solutions are prepared to be within $\pm 2\%$ of the stated concentration, unless otherwise stated on the COA. Cumulative uncertainty from all formulation preparation steps such as weighings and dilutions are presented to help analysts determine the accuracy of their own measurements. Uncertainty for mixtures will of necessity have higher uncertainties than those for individual solutions.

Chemical Purities

- Chemical purities of unlabeled/native standards are 97-99+% unless otherwise specified.
- Chemical purities of labeled standards are 95-99+% unless otherwise specified.

Isotopic Enrichment

Isotopic enrichments specifications for labeled standards are 98-99% for ^{13}C -labeled atoms, 97-98% for deuterium-labeled atoms, and 96% for ^{37}Cl -labeled atoms, unless otherwise specified (refer to product CoA for actual isotopic enrichment). Other isotopic enrichments such as ^{15}N and ^{18}O are on a case-by-case basis.

Solvents Used in Quantitative Solutions

For many years, virtually all of CIL's standards were prepared in nonane or isoctane, solvents that CIL buys in bulk as spectrophotometric grade, and then distills even further to make suitable for ultra-trace analysis.

In recent years, as a larger proportion of standards offered by CIL are more polar, and often analyzed in polar media, many standards are now available in polar solvents such

as MTBE, acetonitrile, methanol, and even water. Many of these standards are unstable in certain solvents, so CIL carefully chooses solvents that are optimal in terms of solubility and stability. Even so, some of these classes of compounds are only stable for a few months to a couple years, and special care should be taken to ensure that they are stored properly. In some cases, co-solvents are required to enable better solubility.

Storage

When standards are frozen, it is a good idea to gently warm them to room temperature and vortex before opening. Sonication can cause decomposition of certain compounds, so it should only be used very carefully.

The biggest threats to the integrity of unopened standards are heat, light (especially UV), oxygen, and pH. These threats also affect opened standards, or standards that have been combined with other compounds, but opened standards are also subject to contamination from glassware/containers, pipettes, and ambient air. The integrity of opened standards is also threatened by solvent evaporation, solubility (when frozen), and decomposition when mixed with other solvents.

It is always good practice to re-certify standards when they have been in storage; the safest laboratory practice is to re-certify standards each time they are used.

Statement on Shelf-Life Testing and Expiration Dates

CIL assigns a retest date and/or expiration date for all manufactured products. The assignments procedure is based upon the known stability characteristics of each product.

- CIL assigns expiration dates for unopened ampoules which have been stored according to CIL's storage instructions.
- CIL's expiration date becomes irrelevant once a standard has been opened and combined with other standards, diluted with solvent, or transferred to a new container. The suitability of the standard then becomes completely dependent on the storage conditions (temperature, light, exposure to other analytes and solvents) and the handling of the standard. The user's QA protocols should determine the duration of the standard's use, and the frequency that it is retested.
- Some products are known to be stable indefinitely. For these products, CIL assigns an expiration date of ten years from the date of release from QC. All other products are retested in five years or less, depending on their chemical characteristics. Most CIL products are sold and shipped with at least one year remaining before the retest date, but they may be shipped with less than one year remaining when products have shorter shelf lives.

Environmental Food, Water, and Exposure Products Information (continued)

Additional Information

24-Hour Emergency Response

CIL and its direct subsidiary CIL Isotope Separations, LLC, are registered with Emergency Response CHEMTREC®. In the event of a chemical-transportation emergency, CHEMTREC® provides immediate advice for those at the scene of emergencies, then promptly contacts the shipper of the chemicals for more detailed assistance and appropriate follow-up. CHEMTREC® operates 24 hours a day, seven

days a week to receive emergency calls. In the case of chemical-transportation emergencies, call one of the following numbers:

Continental United States: 1.800.424.9300

Outside of Continental USA: 1.703.527.3887 (this number may be called collect)

CHEMTREC is a registered trademark of American Chemistry Council, Inc.

Packaging Information



Solids

Solids are packaged in amber glass, wide-mouth, screw-cap jars with teflon-lined caps and tape seals. Clear-glass conical vials are used for small quantities.



Quantitative Solutions

While some of CIL's Environmental Contaminant Standards are packaged in vials, most are supplied as quantitative solutions in amber vials to prevent evaporation. Ampoules are stored in cardboard rondos or plastic "clamshells" with outside labels. Follow the storage instructions, and carefully transfer to other packaging or end-use mixtures as appropriate.

Quality Standards

ISO/IEC 17025

Accreditation to ISO/IEC 17025 "General Requirements for the Competence of Testing and Calibration Laboratories" addresses technical competency of the testing laboratory and specifically, a laboratory's ability to produce precise and accurate test and calibration data. While designed to be applicable to laboratories running third-party samples, accurate test results are critical to the proper characterization of reference materials and, therefore, relevant to reference material producers.



ISO Guide 34

Accreditation to ISO Guide 34 "General Requirements for the Competence of Reference Material Producers" is designed to ensure competency in the manufacture of reference materials and assurance that international guidelines are followed in the production and assignment of material property values. Accreditation to ISO Guide 34 involves not only technical competence and good quality management practices but adds verification of critical production management specific to reference material producers.



Regulatory and Labeling Information

GHS Objectives

The objective of GHS (Globally Harmonized System of Classification and Labeling of Chemicals) is to create an internationally recognizable system for HazCom standards, establish a standard format for hazard communication and support the trade of chemicals for international exchange.

Material Safety Data Sheets (MSDS) are now referred to as Safety Data Sheets (SDS). The SDS functions as an MSDS for ISO, EU and ANSI requirements:

- Most comprehensive information source
- More hazards, including environmental hazards, are now included
- Provides advice and safety precautions
- Product focused; not workplace or task specific
- Written and supplied by manufacturer

Chemical manufacturers/importers/exporters are still responsible for providing information about the identities and hazards of chemicals. All employers using chemicals within their operations are still required to have a hazard communication program.

| GHS Hazard Pictograms and Related Hazards Classes | | |
|---|--|---|
| | | |
| Exploding Bomb • Explosives • Self-reactives • Organic peroxides | Corrosion • Skin corrosion/burns • Eye damage • Corrosive to metals | Flame Over Circle • Oxidizing gases • Oxidizing liquids • Oxidizing solids |
| | | |
| Gas Cylinder • Gases under pressure | Environment • Aquatic toxicity | Skull and Crossbones • Acute toxicity (fatal or toxic) |
| | | |
| Exclamation Point • Irritant (eye and skin) • Skin sensitizer • Acute toxicity • Narcotic effects • Respiratory tract irritant • Hazardous to ozone layer (nonmandatory) | Health Hazard • Carcinogen • Mutagenicity • Reproductive toxicity • Respiratory sensitization • Target organ toxicity • Aspiration toxicity | Flame • Flammables • Pyrophorics • Self-heating • Emits flammable gas • Self-reactives • Organic peroxides |

CIL Labels

Our labels include:

- Product name and description
- Health and safety information
- Lot-specific number
- Package size
- Pictograms for hazard recognition
- CAS numbers
- Storage information
- Packaging number
- Catalog number

Dioxin and Furan Individual Standards



Dioxins and furans are organic pollutants that can be found as byproducts in commercial organochloride pesticide formulations, chlorine-bleached pulp and paper products, and incineration of organic material in the presence of chlorine. Dioxins are persistent in the environment, bioaccumulate in humans and animals, and suspected to be teratogenic, mutagenic, and carcinogenic.

Dioxin and Furan Reference Standards Preparation

The preparation of polychlorinated dibenzo-*p*-dioxin (PCDD) and dibenzofuran (PCDF) certified solution standards begins with the total synthesis of each isomer from known, well-characterized intermediates. Cerilliant QC protocol specifies that all materials be tested to determine identity (multiple techniques), isomer specificity, and purity (multiple techniques), prior to acceptance as a raw material. With few exceptions, our specifications require a chemical purity of >98% for native material and chemical purity of >97% for ¹³C material.

Preparation of CIL/Cerilliant-certified solution standards is tightly controlled using a validated process to ensure accuracy and consistency. Our gravimetric approach (both analyte and solvent are added by weight) is performed using high precision five-place, micro and ultra-micro analytical balances and governed by exacting procedures to ensure minimal uncertainty. Balances are fully qualified in their installed state, are calibrated semi-annually with weekly and pre-use verifications performed – all using NIST traceable weights. Various controls are employed during the dispensing process to ensure no evaporation, degradation, or contamination occurs and to ensure homogeneity and consistency of fill volume from ampoule to ampoule.

Fully certified standards are then put through rigorous QC testing to verify concentration accuracy, consistency with previous lots (when available), and comparison to the corresponding native or ¹³C analog. Finally, homogeneity is ensured through testing of samples pulled during the dispensing process using a random stratified sampling plan. The analytical results are detailed in a comprehensive certificate of analysis (COA) containing complete traceability documentation, which is supplied with each product at no additional charge.

An international round-robin study composed of independent government, commercial, and research laboratories analyzed all 17 CIL/Cerilliant 2,3,7,8-containing polychlorinated dibenzo-*p*-dioxin (PCDD) and dibenzofuran (PCDF) individual solution standards in August 1987. The objective of the study was to determine the accuracy of CIL/Cerilliant solution reference standards. The consensus average values for each of these solutions agreed closely with CIL/Cerilliant reported values – in fact, 15 out of the 17 were within 4%.

Unlabeled Chlorodioxin/Furan Standards for Elution Profiling

CIL offers the only set of all 136 tetra-octa chlorinated dioxin and furan congeners. These qualitative standards are available as ~25 ng/mL solutions in nonane and are used primarily for elution profiling and peak identification. Researchers can utilize this full suite of standards to help identify unknown peaks in their chromatograms and to fully assess samples in environmental forensic cases. Homolog group kits are available, as is a suite of all 136 congeners. See pages 24-25 for full product listings.

Brominated Dioxin and Furan Individual Standards

In an effort to counteract the effects of low-solubility common to many of the higher homolog brominated dioxins and furans, CIL and Cerilliant have reformulated many of these compounds in a cosolvent solution of 70% nonane with 30% toluene. And while most of the brominated dioxins and furans are already formulated at just 5 µg/mL, as detection limits continue to decrease with advances in instrumentation and cleanup procedures, CIL now offers these products as 1.2 mL units.

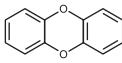
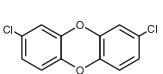
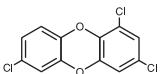
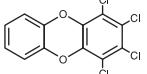
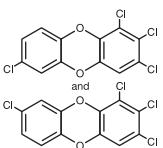
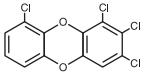
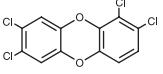
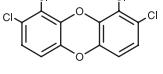
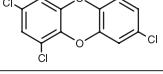
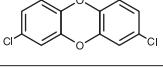
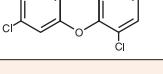
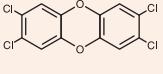
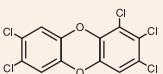
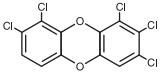
¹³C₁₂-Labeled Chlorodioxin Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------|---|-----------|--|------------|
| CLM-1544-1.2 | Dibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-911 | 1,2,3,4-Tetrachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in nonane 1 ± 0.05 µg/mL in nonane | 1.2 mL |
| ED-911-1 | | | | 1.2 mL |
| ED-4198 | 1,3,6,8-Tetrachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-900 | 2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-4076 | 1,2,3,4,7-Pentachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-955 | 1,2,3,7,8-Pentachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-4077 | 1,2,3,4,6,7-Hexachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-946 | 1,2,3,4,7,8-Hexachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-966 | 1,2,3,6,7,8-Hexachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in 80% nonane/20% toluene | 1.2 mL |
| ED-996 | 1,2,3,7,8,9-Hexachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in 80% nonane/20% toluene | 1.2 mL |
| ED-972 | 1,2,3,4,6,7,8-Heptachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-981 | Octachlorodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) | | 10 ± 0.5 µg/mL in nonane | 4 × 1.2 mL |

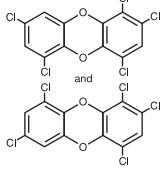
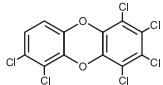
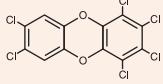
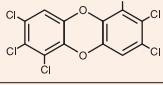
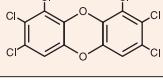
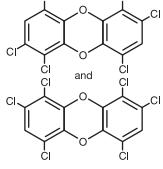
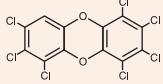
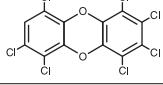
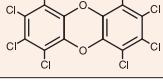
¹³C₆- and ³⁷Cl₄-Labeled Chlorodioxin Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|-------------|---|-----------|------------------------|--------|
| ED-910 | 1,2,3,4-Tetrachlorodibenzo- <i>p</i> -dioxin (¹³ C ₆ , 99%) | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-907 | 2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin (³⁷ Cl ₄ , 96%) | | 50 ± 5 µg/mL in nonane | 1.2 mL |

Unlabeled Chlorodioxin Standards

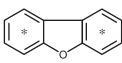
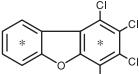
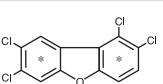
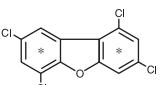
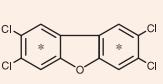
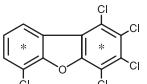
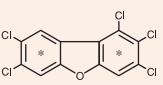
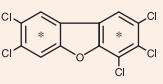
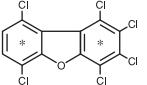
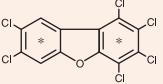
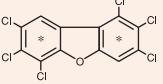
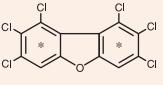
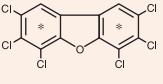
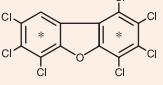
| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------|--|---|-------------------------------|------------|
| ULM-1711-1.2 | Dibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-5568 | 2,8-Dichlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-4090 | 1,3,7-Trichlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-912 | 1,2,3,4-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-905 | 1,2,3,7-Tetrachlorodibenzo- <i>p</i> -dioxin/ 1,2,3,8-Tetrachlorodibenzo- <i>p</i> -dioxin isomer pair |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-948 | 1,2,3,9-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-915 | 1,2,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-916 | 1,2,8,9-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-2518 | 1,3,6,8-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-4061 | 1,3,7,9-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-922 | 1,4,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-901 | 2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane | 4 × 1.2 mL |
| ED-901-A | | | 10 ± 1 µg/mL in methanol | 1.2 mL |
| ED-901-B | | | 50 ± 5 µg/mL in DMSO | 1.2 mL |
| ED-901-C | | | crystalline solid | 1 mg |
| ED-901-D | | | 32 ± 4 pg/µL in DMSO (100 nM) | 0.2 mL |
| ED-950 | 1,2,3,7,8-Pentachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-950-C | | | crystalline solid | 1 mg |
| ED-924 | 1,2,3,8,9-Pentachlorodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |

Unlabeled Chlorodioxin Standards

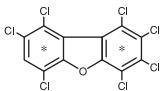
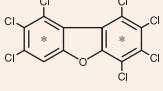
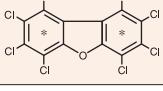
| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------|--|--|---|---------------------|
| ED-927 | 1,2,4,6,8-Pentachlorodibenzo- <i>p</i> -dioxin/ 1,2,4,7,9-Pentachlorodibenzo- <i>p</i> -dioxin isomer pair |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-932 | 1,2,3,4,6,7-Hexachlorodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-961 | 1,2,3,4,7,8-Hexachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-960 ED-960-C | 1,2,3,6,7,8-Hexachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |
| ED-969 ED-969-C | 1,2,3,7,8,9-Hexachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |
| ED-929 | 1,2,4,6,7,9-Hexachlorodibenzo- <i>p</i> -dioxin/ 1,2,4,6,8,9-Hexachlorodibenzo- <i>p</i> -dioxin isomer pair |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-971 | 1,2,3,4,6,7,8-Heptachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| ED-976 | 1,2,3,4,6,7,9-Heptachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| ED-980 ED-980-C | Octachlorodibenzo- <i>p</i> -dioxin |  | 10 ± 0.5 µg/mL in nonane crystalline solid | 4 × 1.2 mL 10 mg |

Other isomers may be available on a special request basis; please inquire.

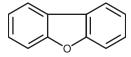
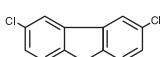
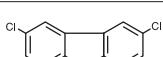
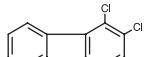
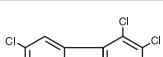
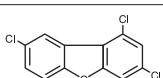
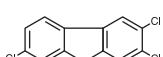
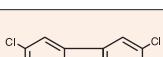
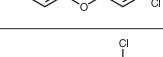
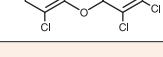
¹³C₁₂-Labeled Chlorofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------|--|--|--------------------------|--------|
| CLM-1561-1.2 | Dibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-920 | 1,2,3,4-Tetrachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-1438 | 1,2,7,8-Tetrachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-5009 | 1,3,6,8-Tetrachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-904 | 2,3,7,8-Tetrachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-5050 | 1,2,3,4,6-Pentachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-952 | 1,2,3,7,8-Pentachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-958 | 2,3,4,7,8-Pentachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-5052 | 1,2,3,4,6,9-Hexachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-963 | 1,2,3,4,7,8-Hexachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-985 | 1,2,3,6,7,8-Hexachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-986 | 1,2,3,7,8,9-Hexachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-987 | 2,3,4,6,7,8-Hexachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-974 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |

¹³C₁₂-Labeled Chlorofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|-------------|--|--|--------------------------|--------|
| EF-5054 | 1,2,3,4,6,8,9-Heptachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-988 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-983 | Octachlorodibenzofuran (¹³ C ₁₂ , 99%) |  | 50 ± 2.5 µg/mL in nonane | 1.2 mL |

Unlabeled Chlorofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------|-----------------------------------|--|---|----------------|
| ULM-1712-1.2 | Dibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-1789 | 2,8-Dichlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-1793 | 2,4,8-Trichlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-921 | 1,2,3,4-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-4030 | 1,2,3,9-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-918 | 1,2,7,8-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-939 | 1,2,8,9-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-944 | 1,3,6,8-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-4031 | 2,3,4,7-Tetrachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-903 EF-903-C | 2,3,7,8-Tetrachlorodibenzofuran |  | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |
| EF-953 EF-953-C | 1,2,3,7,8-Pentachlorodibenzofuran |  | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |
| EF-954 | 1,2,3,8,9-Pentachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-942-50 | 1,3,4,6,8-Pentachlorodibenzofuran |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-956 EF-956-C | 2,3,4,7,8-Pentachlorodibenzofuran |  | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |

Unlabeled Chlorofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------|---------------------------------------|-----------|---|-----------------|
| EF-943-50 | 1,2,3,4,6,8-Hexachlorodibenzofuran | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-964 EF-964-C | 1,2,3,4,7,8-Hexachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 1 mg |
| EF-965 | 1,2,3,4,8,9-Hexachlorodibenzofuran | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EF-962 | 1,2,3,6,7,8-Hexachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-967 | 1,2,3,7,8,9-Hexachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-968 | 2,3,4,6,7,8-Hexachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-973 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-975 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane | 1.2 mL |
| EF-982 EF-982-C | Octachlorodibenzofuran | | 50 ± 2.5 µg/mL in nonane crystalline solid | 1.2 mL 10 mg |

Other isomers may be available on a special request basis; please inquire.

Unlabeled Chlorofuran Standards for Elution Profiling (Qualitative Solutions)

All concentrations are ~25 ng/mL in nonane

| Catalog No. | Compound | Amount |
|-------------|--------------------|--------|
| JR-F01-25 | 1,2,3,4-TetraCDF | 0.2 mL |
| JR-F02-25 | 1,2,3,6-TetraCDF | 0.2 mL |
| JR-F03-25 | 1,2,3,7-TetraCDF | 0.2 mL |
| JR-F04-25 | 1,2,3,8-TetraCDF | 0.2 mL |
| JR-F05-25 | 1,2,3,9-TetraCDF | 0.2 mL |
| JR-F06-25 | 1,2,4,6-TetraCDF | 0.2 mL |
| JR-F07-25 | 1,2,4,7-TetraCDF | 0.2 mL |
| JR-F08-25 | 1,2,4,8-TetraCDF | 0.2 mL |
| JR-F09-25 | 1,2,4,9-TetraCDF | 0.2 mL |
| JR-F10-25 | 1,2,6,7-TetraCDF | 0.2 mL |
| JR-F11-25 | 1,2,6,8-TetraCDF | 0.2 mL |
| JR-F12-25 | 1,2,6,9-TetraCDF | 0.2 mL |
| JR-F13-25 | 1,2,7,8-TetraCDF | 0.2 mL |
| JR-F14-25 | 1,2,7,9-TetraCDF | 0.2 mL |
| JR-F15-25 | 1,2,8,9-TetraCDF | 0.2 mL |
| JR-F16-25 | 1,3,4,6-TetraCDF | 0.2 mL |
| JR-F17-25 | 1,3,4,7-TetraCDF | 0.2 mL |
| JR-F18-25 | 1,3,4,8-TetraCDF | 0.2 mL |
| JR-F19-25 | 1,3,4,9-TetraCDF | 0.2 mL |
| JR-F20-25 | 1,3,6,7-TetraCDF | 0.2 mL |
| JR-F21-25 | 1,3,6,8-TetraCDF | 0.2 mL |
| JR-F22-25 | 1,3,6,9-TetraCDF | 0.2 mL |
| JR-F23-25 | 1,3,7,8-TetraCDF | 0.2 mL |
| JR-F24-25 | 1,3,7,9-TetraCDF | 0.2 mL |
| JR-F25-25 | 1,4,6,7-TetraCDF | 0.2 mL |
| JR-F26-25 | 1,4,6,8-TetraCDF | 0.2 mL |
| JR-F27-25 | 1,4,6,9-TetraCDF | 0.2 mL |
| JR-F28-25 | 1,4,7,8-TetraCDF | 0.2 mL |
| JR-F29-25 | 1,6,7,8-TetraCDF | 0.2 mL |
| JR-F30-25 | 2,3,4,6-TetraCDF | 0.2 mL |
| JR-F31-25 | 2,3,4,7-TetraCDF | 0.2 mL |
| JR-F32-25 | 2,3,4,8-TetraCDF | 0.2 mL |
| JR-F33-25 | 2,3,6,7-TetraCDF | 0.2 mL |
| JR-F34-25 | 2,3,6,8-TetraCDF | 0.2 mL |
| JR-F35-25 | 2,3,7,8-TetraCDF | 0.2 mL |
| JR-F36-25 | 2,4,6,7-TetraCDF | 0.2 mL |
| JR-F37-25 | 2,4,6,8-TetraCDF | 0.2 mL |
| JR-F38-25 | 3,4,6,7-TetraCDF | 0.2 mL |
| JR-F39-25 | 1,2,3,4,6-PentaCDF | 0.2 mL |
| JR-F40-25 | 1,2,3,4,7-PentaCDF | 0.2 mL |
| JR-F41-25 | 1,2,3,4,8-PentaCDF | 0.2 mL |
| JR-F42-25 | 1,2,3,4,9-PentaCDF | 0.2 mL |
| JR-F43-25 | 1,2,3,6,7-PentaCDF | 0.2 mL |
| JR-F44-25 | 1,2,3,6,8-PentaCDF | 0.2 mL |

| Catalog No. | Compound | Amount |
|-------------|-------------------------|--------|
| JR-F45-25 | 1,2,3,6,9-PentaCDF | 0.2 mL |
| JR-F46-25 | 1,2,3,7,8-PentaCDF | 0.2 mL |
| JR-F47-25 | 1,2,3,7,9-PentaCDF | 0.2 mL |
| JR-F48-25 | 1,2,3,8,9-PentaCDF | 0.2 mL |
| JR-F49-25 | 1,2,4,6,7-PentaCDF | 0.2 mL |
| JR-F50-25 | 1,2,4,6,8-PentaCDF | 0.2 mL |
| JR-F51-25 | 1,2,4,6,9-PentaCDF | 0.2 mL |
| JR-F52-25 | 1,2,4,7,8-PentaCDF | 0.2 mL |
| JR-F53-25 | 1,2,4,7,9-PentaCDF | 0.2 mL |
| JR-F54-25 | 1,2,4,8,9-PentaCDF | 0.2 mL |
| JR-F55-25 | 1,2,6,7,8-PentaCDF | 0.2 mL |
| JR-F56-25 | 1,2,6,7,9-PentaCDF | 0.2 mL |
| JR-F57-25 | 1,3,4,6,7-PentaCDF | 0.2 mL |
| JR-F58-25 | 1,3,4,6,8-PentaCDF | 0.2 mL |
| JR-F59-25 | 1,3,4,6,9-PentaCDF | 0.2 mL |
| JR-F60-25 | 1,3,4,7,8-PentaCDF | 0.2 mL |
| JR-F61-25 | 1,3,4,7,9-PentaCDF | 0.2 mL |
| JR-F62-25 | 1,3,6,7,8-PentaCDF | 0.2 mL |
| JR-F63-25 | 1,4,6,7,8-PentaCDF | 0.2 mL |
| JR-F64-25 | 2,3,4,6,7-PentaCDF | 0.2 mL |
| JR-F65-25 | 2,3,4,6,8-PentaCDF | 0.2 mL |
| JR-F66-25 | 2,3,4,7,8-PentaCDF | 0.2 mL |
| JR-F67-25 | 1,2,3,4,6,7-HexaCDF | 0.2 mL |
| JR-F68-25 | 1,2,3,4,6,8-HexaCDF | 0.2 mL |
| JR-F69-25 | 1,2,3,4,6,9-HexaCDF | 0.2 mL |
| JR-F70-25 | 1,2,3,4,7,8-HexaCDF | 0.2 mL |
| JR-F71-25 | 1,2,3,4,7,9-HexaCDF | 0.2 mL |
| JR-F72-25 | 1,2,3,4,8,9-HexaCDF | 0.2 mL |
| JR-F73-25 | 1,2,3,6,7,8-HexaCDF | 0.2 mL |
| JR-F74-25 | 1,2,3,6,7,9-HexaCDF | 0.2 mL |
| JR-F75-25 | 1,2,3,6,8,9-HexaCDF | 0.2 mL |
| JR-F76-25 | 1,2,3,7,8,9-HexaCDF | 0.2 mL |
| JR-F77-25 | 1,2,4,6,7,8-HexaCDF | 0.2 mL |
| JR-F78-25 | 1,2,4,6,7,9-HexaCDF | 0.2 mL |
| JR-F79-25 | 1,2,4,6,8,9-HexaCDF | 0.2 mL |
| JR-F80-25 | 1,3,4,6,7,8-HexaCDF | 0.2 mL |
| JR-F81-25 | 1,3,4,6,7,9-HexaCDF | 0.2 mL |
| JR-F82-25 | 2,3,4,6,7,8-HexaCDF | 0.2 mL |
| JR-F83-25 | 1,2,3,4,6,7,8-HeptaCDF | 0.2 mL |
| JR-F84-25 | 1,2,3,4,6,7,9-HeptaCDF | 0.2 mL |
| JR-F85-25 | 1,2,3,4,6,8,9-HeptaCDF | 0.2 mL |
| JR-F86-25 | 1,2,3,4,7,8,9-HeptaCDF | 0.2 mL |
| JR-F87-25 | 1,2,3,4,6,7,8,9-OctaCDF | 0.2 mL |

Unlabeled Chlorodioxin Standards for Elution Profiling (Qualitative Solutions)

All concentrations are ~25 ng/mL in nonane

| Catalog No. | Compound | Amount |
|-------------|--------------------|--------|
| JR-D01-25 | 1,2,3,4-TetraCDD | 0.2 mL |
| JR-D02-25 | 1,2,3,6-TetraCDD | 0.2 mL |
| JR-D03-25 | 1,2,3,7-TetraCDD | 0.2 mL |
| JR-D04-25 | 1,2,3,8-TetraCDD | 0.2 mL |
| JR-D05-25 | 1,2,3,9-TetraCDD | 0.2 mL |
| JR-D06-25 | 1,2,4,6-TetraCDD | 0.2 mL |
| JR-D07-25 | 1,2,4,7-TetraCDD | 0.2 mL |
| JR-D08-25 | 1,2,4,8-TetraCDD | 0.2 mL |
| JR-D09-25 | 1,2,4,9-TetraCDD | 0.2 mL |
| JR-D10-25 | 1,2,6,7-TetraCDD | 0.2 mL |
| JR-D11-25 | 1,2,6,8-TetraCDD | 0.2 mL |
| JR-D12-25 | 1,2,6,9-TetraCDD | 0.2 mL |
| JR-D13-25 | 1,2,7,8-TetraCDD | 0.2 mL |
| JR-D14-25 | 1,2,7,9-TetraCDD | 0.2 mL |
| JR-D15-25 | 1,2,8,9-TetraCDD | 0.2 mL |
| JR-D16-25 | 1,3,6,8-TetraCDD | 0.2 mL |
| JR-D17-25 | 1,3,6,9-TetraCDD | 0.2 mL |
| JR-D18-25 | 1,3,7,8-TetraCDD | 0.2 mL |
| JR-D19-25 | 1,3,7,9-TetraCDD | 0.2 mL |
| JR-D20-25 | 1,4,6,9-TetraCDD | 0.2 mL |
| JR-D21-25 | 1,4,7,8-TetraCDD | 0.2 mL |
| JR-D22-25 | 2,3,7,8-TetraCDD | 0.2 mL |
| JR-D23-25 | 1,2,3,4,6-PentaCDD | 0.2 mL |
| JR-D24-25 | 1,2,3,4,7-PentaCDD | 0.2 mL |
| JR-D25-25 | 1,2,3,6,7-PentaCDD | 0.2 mL |

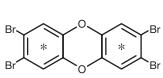
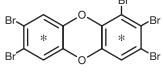
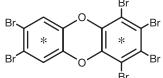
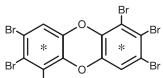
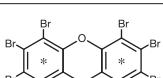
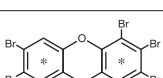
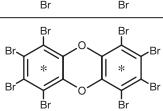
| Catalog No. | Compound | Amount |
|-------------|-------------------------|--------|
| JR-D26-25 | 1,2,3,6,8-PentaCDD | 0.2 mL |
| JR-D27-25 | 1,2,3,6,9-PentaCDD | 0.2 mL |
| JR-D28-25 | 1,2,3,7,8-PentaCDD | 0.2 mL |
| JR-D29-25 | 1,2,3,7,9-PentaCDD | 0.2 mL |
| JR-D30-25 | 1,2,3,8,9-PentaCDD | 0.2 mL |
| JR-D31-25 | 1,2,4,6,7-PentaCDD | 0.2 mL |
| JR-D32-25 | 1,2,4,6,8-PentaCDD | 0.2 mL |
| JR-D33-25 | 1,2,4,6,9-PentaCDD | 0.2 mL |
| JR-D34-25 | 1,2,4,7,8-PentaCDD | 0.2 mL |
| JR-D35-25 | 1,2,4,7,9-PentaCDD | 0.2 mL |
| JR-D36-25 | 1,2,4,8,9-PentaCDD | 0.2 mL |
| JR-D37-25 | 1,2,3,4,6,7-HexaCDD | 0.2 mL |
| JR-D38-25 | 1,2,3,4,6,8-HexaCDD | 0.2 mL |
| JR-D39-25 | 1,2,3,4,6,9-HexaCDD | 0.2 mL |
| JR-D40-25 | 1,2,3,4,7,8-HexaCDD | 0.2 mL |
| JR-D41-25 | 1,2,3,6,7,8-HexaCDD | 0.2 mL |
| JR-D42-25 | 1,2,3,6,7,9-HexaCDD | 0.2 mL |
| JR-D43-25 | 1,2,3,6,8,9-HexaCDD | 0.2 mL |
| JR-D44-25 | 1,2,3,7,8,9-HexaCDD | 0.2 mL |
| JR-D45-25 | 1,2,4,6,7,9-HexaCDD | 0.2 mL |
| JR-D46-25 | 1,2,4,6,8,9-HexaCDD | 0.2 mL |
| JR-D47-25 | 1,2,3,4,6,7,8-HeptaCDD | 0.2 mL |
| JR-D48-25 | 1,2,3,4,6,7,9-HeptaCDD | 0.2 mL |
| JR-D49-25 | 1,2,3,4,6,7,8,9-OctaCDD | 0.2 mL |

Unlabeled Chlorodioxin and Chlorofuran Standard Kits for Elution Profiling

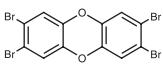
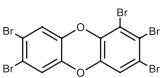
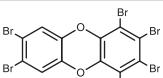
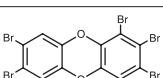
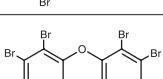
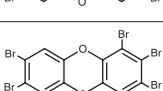
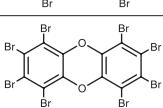
For convenience, CIL has bundled these standards by level of chlorination. Kits are available for tetra-through-hexa dioxins and tetra-through-hepta furans. A comprehensive kit containing all available standards is also available.

| Catalog No. | Description | Contains | Amount |
|---------------|--|-----------------------|--------|
| JR-TCDD-KIT | Comprehensive Tetrachlorodibenzo-p-dioxin Column Defining Kit | JR-D01-25 – JR-D22-25 | 1 Kit |
| JR-TCDF-KIT | Comprehensive Tetrachlorodibenzofuran Column Defining Kit | JR-F01-25 – JR-F38-25 | 1 Kit |
| JR-PECDD-KIT | Comprehensive Pentachlorodibenzo-p-dioxin Column Defining Kit | JR-D23-25 – JR-D36-25 | 1 Kit |
| JR-PECDF-KIT | Comprehensive Pentachlorodibenzofuran Column Defining Kit | JR-F39-25 – JR-F66-25 | 1 Kit |
| JR-HXCDD-KIT | Comprehensive Hexachlorodibenzo-p-dioxin Column Defining Kit | JR-D37-25 – JR-D46-25 | 1 Kit |
| JR-HXCF-KIT | Comprehensive Hexachlorodibenzofuran Column Defining Kit | JR-F67-25 – JR-F82-25 | 1 Kit |
| JR-HPCDF-KIT | Comprehensive Heptachlorodibenzofuran Column Defining Kit | JR-F83-25 – JR-F86-25 | 1 Kit |
| JR-PCDD/F-KIT | Comprehensive Polychlorinated Dioxin and Furan Column Defining Kit (Includes all 136 "JR" dioxin and furan congeners) | | 1 Kit |

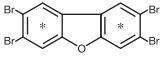
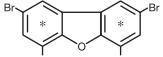
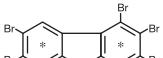
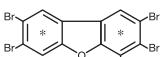
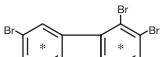
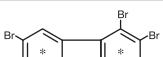
¹³C₁₂-Labeled Bromodioxin Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------------|--|--|---|--------|
| ED-1440-1.2 | 2,3,7,8-Tetrabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-1450-1.2 | 1,2,3,7,8-Pentabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| NEW ED-2534-A-1.2 | 1,2,3,4,7,8-Hexabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| NEW ED-5237-A-1.2 | 1,2,3,6,7,8-Hexabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| NEW ED-5238-A-1.2 | 1,2,3,7,8,9-Hexabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| ED-5357-1.2 | 1,2,3,4,6,7,8-Heptabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| ED-5089-1.2 | Octabromodibenzo- <i>p</i> -dioxin (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |

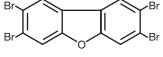
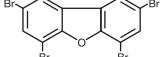
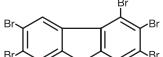
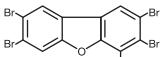
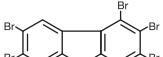
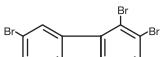
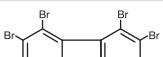
Unlabeled Bromodioxin Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------------|---|--|---|--------|
| ED-1441-1.2 | 2,3,7,8-Tetrabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| ED-1451-1.2 | 1,2,3,7,8-Pentabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| NEW ED-1462-A-1.2 | 1,2,3,4,7,8-Hexabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| ED-1465-1.2 | 1,2,3,6,7,8-Hexabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| NEW ED-1466-A-1.2 | 1,2,3,7,8,9-Hexabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| ED-5356-1.2 | 1,2,3,4,6,7,8-Heptabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| ED-1481-1.2 | Octabromodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in toluene | 1.2 mL |

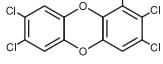
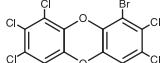
¹³C₁₂-Labeled Bromofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|-------------|---|--|---|--------|
| EF-1442-1.2 | 2,3,7,8-Tetrabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-5082-1.2 | 2,4,6,8-Tetrabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1452-1.2 | 1,2,3,7,8-Pentabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1454-1.2 | 2,3,4,7,8-Pentabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1463-1.2 | 1,2,3,4,7,8-Hexabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-5259-1.2 | 1,2,3,4,6,7,8-Heptabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| EF-5266-1.2 | Octabromodibenzofuran (¹³ C ₁₂ , 99%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |

Unlabeled Bromofuran Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|------------------------|---|--|---|--------|
| EF-1443-1.2 | 2,3,7,8-Tetrabromodibenzofuran |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| NEW EF-5081-1.2 | 2,4,6,8-Tetrabromodibenzofuran |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1453-1.2 | 1,2,3,7,8-Pentabromodibenzofuran |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1455-1.2 | 2,3,4,7,8-Pentabromodibenzofuran |  | 5 ± 0.5 µg/mL in nonane | 1.2 mL |
| EF-1464-1.2 | 1,2,3,4,7,8-Hexabromodibenzofuran |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| EF-1486-1.2 | 1,2,3,4,6,7,8-Heptabromodibenzofuran (CP 96%) |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |
| EF-5263-1.2 | Octabromodibenzofuran |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |

Unlabeled Mixed Bromo/Chlorodioxin Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------|---|--|---|--------|
| NEW EBC-2501 | 1-Bromo-2,3,7,8-tetrachlorodibenzo- <i>p</i> -dioxin |  | 50 ± 0.5 µg/mL in <i>n</i> -nonane | 1.2 mL |
| NEW | EBC-2507-A-1.2 1-Bromo-2,3,4,6,7,8,9-heptachlorodibenzo- <i>p</i> -dioxin |  | 5 ± 0.5 µg/mL in 70% nonane/30% toluene | 1.2 mL |

Dioxin and Furan Method Standards, Standard Mixtures and Reference Materials



Following the unprecedented release of dioxins and furans into the environment during the Seveso, Italy, industrial accident in 1976, regulatory agencies around the world initiated strict testing programs to monitor these toxic compounds. CIL and Cerilliant collaborated with the US EPA to prepare the first isotope dilution mass spectrometry (IDMS) standard mixtures for the analysis of PCDDs and PCDFs.

US EPA, JIS, and CEN Dioxin and Furan Method Standard Mixtures

In 1990 CIL/Cerilliant (formerly Radian) introduced the first "ready-to-use" standard mixtures for US EPA Method 1613 "High Resolution GC/MS Method for the Determination of Tetra-Octa Chlorinated Dioxins and Furans." With the effectiveness and popularity of these preformulated mixtures, CIL/Cerilliant next developed "ready-to-use" standards for EPA Method 8280 for low-resolution GC/MS analysis of dioxins and furans. Today CIL/Cerilliant offer convenient dioxin and furan standard mixtures for EPA Methods 23 and 8290, as well as the Japanese Industrial Standards methods JIS-K0311 and K0312, and the European Community method EN-1948. Copies of these methods are available upon request.

Dioxin and Furan Plus PCB Standard Mixtures

CIL/Cerilliant have developed several mixtures that include the 2,3,7,8-containing dioxin and furan congeners, as well as the "toxic" PCB congeners. These mixtures have full calibration series and matching spiking solutions, so analysts can test these two commonly combined groups without having to manipulate several different standard sets.

Dioxin and Furan "Starter Kits"

CIL recognizes that starting up a dioxin and furan testing laboratory can be a daunting and expensive process. To assist laboratories that are just getting started, CIL offers "starter kits" to help procuring the standards needed for dioxin and furan testing much easier. These kits include all required standard mixtures, including calibration and spiking cocktails, for specific testing methods. There is no guesswork required for materials that will be needed to start the process, and a kit will typically test between 50-200 samples. Once the method is established, it will be clear which standards will be needed sooner and those that can be reordered later. Most standards last several years, so there is no concern for shelf life. In addition, the cost of the kits has been reduced significantly (when compared to purchasing all mixtures separately), easing the financial burden of setting up a new lab!

Reference Materials

In 2006, CIL completed an international interlaboratory study for the determination of many environmental pollutants in its three fish reference materials, as well as two reference materials for soil and sediment. In 2007, CIL conducted another interlaboratory study, this time evaluating dioxins, furans, and PCBs in a new fly ash reference material. In 2010, CIL launched yet another interlaboratory study to develop consensus values for priority pollutants in cod liver oil reference material.

Non-2,3,7,8-Containing Standard Mixtures

With the development of several ¹³C-labeled "non-2,3,7,8" furan standards, CIL/Cerilliant can offer standard mixtures that contain the traditional 17 "2,3,7,8-containing" standards, as well as ¹³C-labeled "non-2,3,7,8-containing" congeners. These standard mixtures allow researchers to use all 17 ¹³C-labeled 2,3,7,8-containing standards as internal standards, while utilizing the labeled "non-2,3,7,8-containing" congeners as recovery/injection or cleanup standards.

Two-Column Dioxin and Furan Standard Mixtures

Two-column dioxin and furan standard mixtures are combination mixtures used to confirm dioxins and furans and PCBs using only two columns. These standards combine the benefits of both the "dioxin and furan plus PCB" mixtures and the "non-2,3,7,8-containing" mixtures.

Expanded PBDD/F Standards and Standard Mixtures

Polybrominated dioxins and furans (PBDD/F) can be found at trace levels in technical brominated flame-retardant products, and may also be formed from combustion of these materials in the presence of organic compounds. The biological effects of PBDD/Fs are similar to those of their chlorinated analogs, which have been regulated for many years. CIL offers a comprehensive set of labeled and unlabeled standards for PBDD/F analysis, including calibration series and corresponding spiking solutions containing tetra-octabromo congeners.

US EPA Method 1613 Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|--|-----------------------------|
| EDF-9999 | Method 1613 Calibration Solutions [CS1-CS5] | Set of 5 × 0.2 mL in nonane |
| * EDF-9999-0.1 | Method 1613 Calibration Solution [CS0.1] | 0.2 mL in nonane |
| * EDF-9999-0.2 | Method 1613 Calibration Solution [CS0.2] | 0.2 mL in nonane |
| * EDF-9999-0.5 | Method 1613 Calibration Solution [CS0.5] | 0.2 mL in nonane |
| EDF-9999-1 | Method 1613 Calibration Solution [CS1] | 0.2 mL in nonane |
| EDF-9999-2 | Method 1613 Calibration Solution [CS2] | 0.2 mL in nonane |
| EDF-9999-3 | Method 1613 Daily Calibration Check Standard [CS3] | 0.2 mL in nonane |
| EDF-9999-3-4 | Method 1613 Daily Calibration Check Standard [CS3] | Set of 4 × 0.2 mL in nonane |
| EDF-9999-4 | Method 1613 Calibration Solution [CS4] | 0.2 mL in nonane |
| EDF-9999-5 | Method 1613 Calibration Solution [CS5] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | *CS0.1 | *CS0.2 | *CS0.5 | CS1 | CS2 | CS3 | CS4 | CS5 |
|---|--------|--------|--------|-----|-----|-----|-----|------|
| 2,3,7,8-TetraCDD | 0.05 | 0.1 | 0.25 | 0.5 | 2.0 | 10 | 40 | 200 |
| 2,3,7,8-TetraCDF | 0.05 | 0.1 | 0.25 | 0.5 | 2.0 | 10 | 40 | 200 |
| 1,2,3,7,8-PentaCDD | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,7,8-PentaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,4,7,8-PentaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8-HexaCDD | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,6,7,8-HexaCDD | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,7,8,9-HexaCDD | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8-HexaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,6,7,8-HexaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,7,8,9-HexaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 2,3,4,6,7,8-HexaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.25 | 0.5 | 1.25 | 2.5 | 10 | 50 | 200 | 1000 |
| OctaCDD | 0.5 | 1.0 | 2.50 | 5.0 | 20 | 100 | 400 | 2000 |
| OctaCDF | 0.5 | 1.0 | 2.50 | 5.0 | 20 | 100 | 400 | 2000 |
| Labeled | | | | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| OctaCDD (¹³ C ₁₂ , 99%) | 200 | 200 | 200 | 200 | 200 | 200 | 200 | 200 |
| Cleanup | | | | | | | | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 0.05 | 0.1 | 0.25 | 0.5 | 2.0 | 10 | 40 | 200 |
| Internal | | | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |

*NOTE: CS0.1, CS0.2, and CS0.5 are optional extensions of the Method 1613 Calibration Curve to extend the MDL and are not required by the method.

US EPA Method 1613 Standard Mixtures

| Catalog No. | Compound | Amount | | | | |
|---|---|-----------------------------|-----|-----|-----|-----|
| EDF-9999-A | Method 1613 Calibration Solutions (1/10 concentration) [CS1-CS5] | Set of 5 x 0.2 mL in nonane | | | | |
| EDF-9999-A-3 | Method 1613 Calibration Check Standard (1/10 concentration) [CS3] | 0.2 mL in nonane | | | | |
| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | |
| Unlabeled | | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,3,7,8-TetraCDD | 0.05 | 0.2 | 1 | 4 | 20 | |
| 2,3,7,8-TetraCDF | 0.05 | 0.2 | 1 | 4 | 20 | |
| 1,2,3,7,8-PentaCDD | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,7,8-PentaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 2,3,4,7,8-PentaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,4,7,8-HexaCDD | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,6,7,8-HexaCDD | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,7,8,9-HexaCDD | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,4,7,8-HexaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,6,7,8-HexaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,7,8,9-HexaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 2,3,4,6,7,8-HexaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,4,6,7,8-HeptaCDD | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,4,6,7,8-HeptaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| 1,2,3,4,7,8,9-HeptaCDF | 0.25 | 1 | 5 | 20 | 100 | |
| OctaCDD | 0.5 | 2 | 10 | 40 | 200 | |
| OctaCDF | 0.5 | 2 | 10 | 40 | 200 | |
| Labeled | | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | |
| Cleanup | | | | | | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 99%) | 0.05 | 0.2 | 1 | 4 | 20 | |
| Internal | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | |

EDF-9999-A is a set of calibration solutions with both labeled and unlabeled compounds at 1/10 the concentration of the corresponding calibration solution in EDF-9999.

| | | |
|--|--|-------|
| EDF-1613-KIT | Method 1613 "Starter Kit" | 1 Kit |
| <i>Contains one each of the following items:</i> | | |
| EDF-9999 | Method 1613 Calibration Solutions [CS1-CS5] | |
| EDF-8999(2X) | Method 1613 Labeled Compound Stock Solution | |
| EDF-5999 | Method 1613 Internal Standard Spiking Solution | |
| EDF-7999 | Method 1613 Precision and Recovery Standard Solution | |
| EDF-6999 | Method 1613 Cleanup Standard | |

US EPA Method 1613 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|---|-----------------------------|
| EDF-8999 | Method 1613 Labeled Compound Stock Solution | 500 µL in nonane |
| EDF-8999-4 | Method 1613 Labeled Compound Stock Solution | Set of 4 × 500 µL in nonane |
| NEW EDF-8999-5ML | Method 1613 Labeled Compound Stock Solution | 5 mL in nonane |

| Labeled | (ng/mL) |
|---|---------|
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 100 |
| OctaCDD (¹³ C ₁₂ , 99%) | 200 |

| | | |
|--------------|--|------------------|
| EDF-6999 | Method 1613 Cleanup Standard | 7.5 mL in nonane |
| EDF-6999-10X | Method 1613 Cleanup Standard (10X concentration) | 20 mL in nonane |

| Labeled | EDF-6999 (ng/mL) | EDF-6999-10X (ng/mL) |
|---|---------------------|-------------------------|
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 0.8 | 8 |

| | | |
|---------------------------|--|------------------|
| EDF-5999 | Method 1613 Internal Standard Spiking Solution | 0.5 mL in nonane |
| NEW EDF-5999-2.5ML | Method 1613 Internal Standard Spiking Solution | 2.5 mL in nonane |

| Labeled | (ng/mL) |
|--|---------|
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 200 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 200 |

| | | |
|--------------|--|------------------|
| EDF-7999 | Method 1613 Precision and Recovery Standard Solution | 200 µL in nonane |
| EDF-7999-10X | Method 1613 Precision and Recovery Standard Solution (10X concentration) | 1.2 mL in nonane |

| Unlabeled | EDF-7999 (ng/mL) | EDF-7999-10X (ng/mL) |
|------------------------|---------------------|-------------------------|
| 2,3,7,8-TetraCDD | 40 | 400 |
| 2,3,7,8-TetraCDF | 40 | 400 |
| 1,2,3,7,8-PentaCDD | 200 | 2000 |
| 1,2,3,7,8-PentaCDF | 200 | 2000 |
| 2,3,4,7,8-PentaCDF | 200 | 2000 |
| 1,2,3,4,7,8-HexaCDD | 200 | 2000 |
| 1,2,3,6,7,8-HexaCDD | 200 | 2000 |
| 1,2,3,7,8,9-HexaCDD | 200 | 2000 |
| 1,2,3,4,7,8-HexaCDF | 200 | 2000 |
| 1,2,3,6,7,8-HexaCDF | 200 | 2000 |
| 1,2,3,7,8,9-HexaCDF | 200 | 2000 |
| 2,3,4,6,7,8-HexaCDF | 200 | 2000 |
| 1,2,3,4,6,7,8-HeptaCDD | 200 | 2000 |
| 1,2,3,4,6,7,8-HeptaCDF | 200 | 2000 |
| 1,2,3,4,7,8,9-HeptaCDF | 200 | 2000 |
| OctaCDD | 400 | 4000 |
| OctaCDF | 400 | 4000 |

US EPA Method 1613 Standard Mixtures

| Catalog No. | Compound | Amount |
|---|---|------------------|
| EDF-4141 | Method 1613 Daily Calibration Plus Window Definer and Isomer Specificity Solution | 200 µL in nonane |
| Daily Calibration | | |
| 2,3,7,8-TetraCDD | (ng/mL) | |
| 2,3,7,8-TetraCDF | 10 | |
| 1,2,3,7,8-PentaCDD | 10 | |
| 1,2,3,7,8-PentaCDF | 50 | |
| 2,3,4,7,8-PentaCDF | 50 | |
| 1,2,3,4,7,8-HexaCDD | 50 | |
| 1,2,3,6,7,8-HexaCDD | 50 | |
| 1,2,3,7,8,9-HexaCDD | 50 | |
| 1,2,3,4,7,8-HexaCDF | 50 | |
| 1,2,3,6,7,8-HexaCDF | 50 | |
| 1,2,3,7,8,9-HexaCDF | 50 | |
| 2,3,4,6,7,8-HexaCDF | 50 | |
| 1,2,3,4,6,7,8-HeptaCDD (WD) | 50 | |
| 1,2,3,4,6,7,8-HeptaCDF (WD) | 50 | |
| 1,2,3,4,7,8,9-HeptaCDF (WD) | 50 | |
| OctaCDD | 100 | |
| OctaCDF | 100 | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 10 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 200 | |
| Window Defining | | |
| 1,3,6,8-TetraCDD | (ng/mL) | |
| 1,2,8,9-TetraCDD | 10 | |
| 1,3,6,8-TetraCDF | 10 | |
| 1,2,8,9-TetraCDF | 10 | |
| 1,2,4,6,8/1,2,4,7,9-PentaCDD | 50 | |
| 1,2,3,8,9-PentaCDD | 50 | |
| 1,3,4,6,8-PentaCDF | 50 | |
| 1,2,3,8,9-PentaCDF | 50 | |
| 1,2,4,6,7,9/1,2,4,6,8,9-HexaCDD | 50 | |
| 1,2,3,4,6,8-HexaCDF | 50 | |
| 1,2,3,4,8,9-HexaCDF | 50 | |
| 1,2,3,4,6,7,9-HeptaCDD | 50 | |

NOTE: 1,2,3,4,6,7-HexaCDD (last eluting HexaCDD) not included due to interference with 1,2,3,7,8,9-HexaCDD.

| TetraCDD Isomer Specificity | (ng/mL) |
|-----------------------------|---------|
| 1,2,3,4-TetraCDD | 10 |
| 1,2,3,7/1,2,3,8-TetraCDD | 10 |
| 1,2,3,9-TetraCDD | 10 |

This standard allows three functions:

- Daily MS instrument calibration verification
- Daily TetraCDD column resolution
- Daily window definition

2,3,7,8-TetraCDD Only Standard Mixture

| Catalog No. | Compound | Amount |
|---|--|------------------|
| NEW ED-4159-CS1 | Calibration Curve, 2,3,7,8 CS1 for 2,3,7,8-TCDD Only | 1.2 mL in nonane |
| Unlabeled | | |
| 2,3,7,8-TetraCDD | (ng/mL) | |
| 2,3,7,8-TetraCDD | 0.01 | |
| Labeled | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 25 | |

Performance Evaluation Reference Materials

Fish Tissue, Soil, and Sediment Reference Materials

CIL first introduced fish tissue reference material (RM) standards in the early 1990s to provide analytical laboratories testing dioxin, furan, and PCB congeners with much needed performance evaluation materials. At that time, these RMs were quite uncommon in the field of dioxin and PCB analysis, and they became very popular for their ability to help laboratories internally assess their quality-assurance processes.

Reference materials are widely used in analytical chemistry to provide immediate feedback about the quality and confidence of a laboratory's testing system. RMs can be used to demonstrate to quality auditors that a laboratory has control of their QA process and can show long-term trending by using the same RM over time.

The original "clean," "fortified," and "naturally contaminated" fish RMs were subsequently joined by natural matrix soil and sediment, cod liver oil, and even chicken egg RMs. CIL RMs often have consensus values for many other organic contaminants, including dioxins, furans, and PCBs, as well as PAHs, pesticides, flame retardants, and others. All of CIL's RM standards are run through an international interlaboratory study to determine consensus values with uncertainties.

| Catalog No. | Compound | Amount |
|---------------------|--|----------|
| EDF-5183 | Natural Matrix Reference Material (Soil) | 10 g |
| EDF-5184 | Heavily Contaminated Sediment Reference Material | 10 g |
| EDF-2524 | Clean Natural Matrix Reference Material (Fish) | 10 g |
| EDF-2525 | Contaminated Natural Matrix Reference Material (Fish) | 10 g |
| EDF-2526 | Fortified Natural Matrix Reference Material (Fish) | 10 g |
| EDF-4023 | Set of 3 Fish (1 each of EDF-2524, EDF-2525, EDF-2526) | 3 × 10 g |
| EDF-5462 | Fortified Cod Liver Oil Reference Material | 10 g |
| EDF-5463 | Cod Liver Oil Reference Material | 10 g |
| NEW EDF-5491 | Freeze-Dried Eggs Reference Material | 6 g |

Perfluorokerosene (PFK)

| Catalog No. | Compound | Amount |
|-------------------------|---|--------|
| NEW PFK-HIGH-0.1 | Perfluorokerosene, High-Boiling Range (unlabeled) | 0.1 g |
| NEW PFK-HIGH-0.5 | Perfluorokerosene, High-Boiling Range (unlabeled) | 0.5 g |
| NEW PFK-HIGH-1 | Perfluorokerosene, High-Boiling Range (unlabeled) | 1 g |
| NEW PFK-LOW-0.25 | Perfluorokerosene, Low-Boiling Range (unlabeled) | 0.25 g |
| NEW PFK-LOW-1 | Perfluorokerosene, Low-Boiling Range (unlabeled) | 1 g |

US EPA Method 23 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|-----------------------------|
| EDF-4052 | Method 23 Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EDF-4052-1 | Method 23 Calibration Solution [CS1] | 0.2 mL in nonane |
| EDF-4052-2 | Method 23 Calibration Solution [CS2] | 0.2 mL in nonane |
| EDF-4052-3 | Method 23 Daily Calibration Check Standard [CS3] | 0.2 mL in nonane |
| EDF-4052-4 | Method 23 Calibration Solution [CS4] | 0.2 mL in nonane |
| EDF-4052-5 | Method 23 Calibration Solution [CS5] | 0.2 mL in nonane |

| <i>All concentrations are in pg/µL (ppb)</i> | | | | | |
|---|-----|-----|-----|-----|------|
| | CS1 | CS2 | CS3 | CS4 | CS5 |
| Unlabeled | | | | | |
| 2,3,7,8-TetraCDD | 0.5 | 1 | 5 | 50 | 100 |
| 2,3,7,8-TetraCDF | 0.5 | 1 | 5 | 50 | 100 |
| 1,2,3,7,8-PentaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,7,8-PentaCDF | 2.5 | 5 | 25 | 250 | 500 |
| 2,3,4,7,8-PentaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-HexaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,6,7,8-HexaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,7,8,9-HexaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-HexaCDF | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,6,7,8-HexaCDF | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,7,8,9-HexaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 2,3,4,6,7,8-HexaCDF | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,6,7,8-HeptaCDD | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,6,7,8-HeptaCDF | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8,9-HeptaCDD | 2.5 | 5 | 25 | 250 | 500 |
| OctaCDD | 5.0 | 10 | 50 | 500 | 1000 |
| OctaCDF | 5.0 | 10 | 50 | 500 | 1000 |
| Internal | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| OctaCDD (¹³ C ₁₂ , 99%) | 200 | 200 | 200 | 200 | 200 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| Surrogate | | | | | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 0.5 | 1 | 5 | 50 | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2.5 | 5 | 25 | 250 | 500 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 2.5 | 5 | 25 | 250 | 500 |
| Recovery | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| Alternate Recovery | | | | | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 2.5 | 5 | 25 | 250 | 500 |

| | | | |
|------------|------------|-------------------------|-------|
| NEW | EDF-23-KIT | Method 23 "Starter Kit" | 1 Kit |
|------------|------------|-------------------------|-------|

Contains one each of the following items:

| | |
|----------|--|
| EDF-4052 | Method 23 Calibration Solutions [CS1-CS5] |
| EDF-4053 | Method 23 Internal Standard Stock Solution |
| EDF-4054 | Method 23 Surrogate Standard Stock Solution |
| EDF-4055 | Method 23 Recovery Standard Stock Solution |
| EDF-5189 | Method 23 Alternate Recovery Standard Stock Solution |

US EPA Method 23 Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| EDF-4053 | Method 23 Internal Standard Stock Solution | 1.2 mL in nonane |
| Labeled (pg/µL) | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| <hr/> | | |
| EDF-4054 | Method 23 Surrogate Standard Stock Solution | 1.2 mL in nonane |
| Labeled (pg/µL) | | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 1000 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| <hr/> | | |
| EDF-4055 | Method 23 Recovery Standard Stock Solution | 1.2 mL in nonane |
| Labeled (pg/µL) | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 500 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 500 | |
| <hr/> | | |
| EDF-5189 | Method 23 Alternate Recovery Standard Stock Solution | 1.2 mL in nonane |
| Labeled (pg/µL) | | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | |

US EPA Method 8290 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EDF-5006 | Method 8290 Calibration Solutions [HRCC1-HRCC5] | Set of 5 x 0.2 mL in nonane |
| EDF-5006-1 | Method 8290 Calibration Solution [HRCC1] | 0.2 mL in nonane |
| EDF-5006-2 | Method 8290 Calibration Solution [HRCC2] | 0.2 mL in nonane |
| EDF-5006-3 | Method 8290 Continuing Calibration Check Standard [HRCC3] | 0.2 mL in nonane |
| EDF-5006-4 | Method 8290 Calibration Solution [HRCC4] | 0.2 mL in nonane |
| EDF-5006-5 | Method 8290 Calibration Solution [HRCC5] | 0.2 mL in nonane |

| <i>All concentrations are in pg/µL (ppb)</i> | | | | | |
|---|-------|-------|-------|-------|-------|
| Unlabeled | HRCC1 | HRCC2 | HRCC3 | HRCC4 | HRCC5 |
| 2,3,7,8-TetraCDD | 1.0 | 2.5 | 10 | 50 | 200 |
| 2,3,7,8-TetraCDF | 1.0 | 2.5 | 10 | 50 | 200 |
| 1,2,3,7,8-PentaCDD | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,7,8-PentaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 2,3,4,7,8-PentaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8-HexaCDD | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,6,7,8-HexaCDD | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,7,8,9-HexaCDD | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8-HexaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,6,7,8-HexaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,7,8,9-HexaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 2,3,4,6,7,8-HexaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,6,7,8-HeptaCDD | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,6,7,8-HeptaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| 1,2,3,4,7,8,9-HeptaCDF | 2.5 | 6.25 | 25 | 125 | 500 |
| OctaCDD | 5.0 | 12.5 | 50 | 250 | 1000 |
| OctaCDF | 5.0 | 12.5 | 50 | 250 | 1000 |
| Internal | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 | 50 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 125 | 125 | 125 | 125 | 125 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 125 | 125 | 125 | 125 | 125 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 125 | 125 | 125 | 125 | 125 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 125 | 125 | 125 | 125 | 125 |
| OctaCDD (¹³ C ₁₂ , 99%) | 250 | 250 | 250 | 250 | 250 |
| Recovery | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 | 50 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 125 | 125 | 125 | 125 | 125 |

| | | | |
|------------|--------------|---------------------------|-------|
| NEW | EDF-8290-KIT | Method 8290 "Starter Kit" | 1 Kit |
|------------|--------------|---------------------------|-------|

Contains one each of the following items:

| | |
|----------|---|
| EDF-5006 | Method 8290 Calibration Solutions [HRCC1-HRCC5] |
| EDF-5005 | Method 8290 Sample Fortification Solution |
| EDF-5004 | Method 8290 Recovery Standard Solution |
| EDF-5008 | Method 8290 Matrix Spiking Solution |

US EPA Method 8290 Standard Mixtures

| Catalog No. | Compound | Amount |
|---|---|------------------|
| EDF-5005 | Method 8290 Sample Fortification Solution | 1.2 mL in nonane |
| Labeled | | (pg/µL) |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 100 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 250 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 250 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 250 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 250 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 500 |
| ED-5004 | Method 8290 Recovery Standard Solution | 1.2 mL in nonane |
| Labeled | | (pg/µL) |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | | 100 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 250 |
| EDF-5008 | Method 8290 Matrix Spiking Solution | 1.2 mL in nonane |
| EDF-5008-50 | Method 8290 Matrix Spiking Solution (1:50 dilution) | 0.1 mL in nonane |
| Unlabeled | | |
| | EDF-5008 | EDF-5008-50 |
| | (pg/µL) | (pg/µL) |
| 2,3,7,8-TetraCDD | 100 | 2 |
| 2,3,7,8-TetraCDF | 100 | 2 |
| 1,2,3,7,8-PentaCDD | 250 | 5 |
| 1,2,3,7,8-PentaCDF | 250 | 5 |
| 2,3,4,7,8-PentaCDD | 250 | 5 |
| 1,2,3,4,7,8-HexaCDD | 250 | 5 |
| 1,2,3,4,7,8-HexaCDF | 250 | 5 |
| 1,2,3,6,7,8-HexaCDD | 250 | 5 |
| 1,2,3,6,7,8-HexaCDF | 250 | 5 |
| 1,2,3,7,8,9-HexaCDD | 250 | 5 |
| 1,2,3,7,8,9-HexaCDF | 250 | 5 |
| 2,3,4,6,7,8-HexaCDD | 250 | 5 |
| 1,2,3,4,6,7,8-HeptaCDD | 250 | 5 |
| 1,2,3,4,6,7,8-HeptaCDF | 250 | 5 |
| 1,2,3,4,7,8,9-HeptaCDD | 250 | 5 |
| OctaCDD | 500 | 10 |
| OctaCDF | 500 | 10 |

European Air Method EN-1948 Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------|---|-----------------------------|
| EDF-4947 | EN-1948 Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EDF-4947-CS1 | EN-1948 Calibration Solution [CS1] | 0.2 mL in nonane |
| EDF-4947-CS2 | EN-1948 Calibration Solution [CS2] | 0.2 mL in nonane |
| EDF-4947-CS3 | EN-1948 Calibration Solution [CS3] | 0.2 mL in nonane |
| EDF-4947-CS4 | EN-1948 Calibration Solution [CS4] | 0.2 mL in nonane |
| EDF-4947-CS5 | EN-1948 Calibration Solution [CS5] | 0.2 mL in nonane |

| <i>All concentrations are in pg/µL (ppb)</i> | | | | | |
|---|-----|------|------|------|------|
| Unlabeled | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,3,7,8-TetraCDD | 0.5 | 2.0 | 10.0 | 40.0 | 200 |
| 2,3,7,8-TetraCDF | 0.5 | 2.0 | 10.0 | 40.0 | 200 |
| 1,2,3,7,8-PentaCDD | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,7,8-PentaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 2,3,4,7,8-PentaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,4,7,8-HexaCDD | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,6,7,8-HexaCDD | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,7,8,9-HexaCDD | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,4,7,8-HexaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,6,7,8-HexaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,7,8,9-HexaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 2,3,4,6,7,8-HexaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDD | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| 1,2,3,4,7,8,9-HeptaCDF | 2.5 | 10.0 | 50.0 | 200 | 1000 |
| OctaCDD | 5.0 | 20.0 | 100 | 400 | 2000 |
| OctaCDF | 5.0 | 20.0 | 100 | 400 | 2000 |
| Sampling | | | | | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| Extraction | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| OctaCDD (¹³ C ₁₂ , 99%) | 200 | 200 | 200 | 200 | 200 |
| OctaCDF (¹³ C ₁₂ , 99%) | 200 | 200 | 200 | 200 | 200 |
| Syringe | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |

| | | | |
|------------|--------------|------------------------------------|-------|
| NEW | EDF-1948-KIT | EN-1948 Dioxin/Furan "Starter Kit" | 1 Kit |
|------------|--------------|------------------------------------|-------|

Contains one each of the following items:

| | |
|----------|---|
| EDF-4947 | EN-1948 Calibration Solutions [CS1-CS5] |
| EF-4138 | EN-1948 Sampling Standard Solution |
| EDF-4139 | EN-1948 Extraction Standard Solution |
| ED-4140 | EN-1948 Syringe Standard Solution |
| EDF-4175 | EN-1948 Native Stock Response Factor Solution |

European Air Method EN-1948 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|--------------------|
| EF-4138 | EN-1948 Sampling Standard Solution | 1.2 mL in nonane |
| EF-4138-10 | EN-1948 Sampling Standard Solution | 2 × 5 mL in nonane |
| | Labeled | (pg/µL) |
| | 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 200 |
| EDF-4139 | EN-1948 Extraction Standard Solution | 1.2 mL in nonane |
| EDF-4139-10 | EN-1948 Extraction Standard Solution | 2 × 5 mL in nonane |
| | Labeled | (pg/µL) |
| | 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 100 |
| | 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 100 |
| | 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 100 |
| | 2,3,4,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 200 |
| | 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 200 |
| | OctaCDD (¹³ C ₁₂ , 99%) | 200 |
| | OctaCDF (¹³ C ₁₂ , 99%) | 200 |
| ED-4140 | EN-1948 Syringe Standard Solution | 1.2 mL in nonane |
| | Labeled | (pg/µL) |
| | 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 400 |
| | 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 400 |
| EDF-4175 | EN-1948 Native Stock Response Factor Solution | 0.5 mL in nonane |
| | Unlabeled | (ng/mL) |
| | 2,3,7,8-TetraCDD | 1000 |
| | 2,3,7,8-TetraCDF | 1000 |
| | 1,2,3,7,8-PentaCDD | 1000 |
| | 1,2,3,7,8-PentaCDF | 1000 |
| | 2,3,4,7,8-PentaCDF | 1000 |
| | 1,2,3,4,7,8-HexaCDD | 1000 |
| | 1,2,3,6,7,8-HexaCDD | 1000 |
| | 1,2,3,7,8,9-HexaCDD | 4000 |
| | 1,2,3,4,7,8-HexaCDF | 1000 |
| | 1,2,3,6,7,8-HexaCDF | 1000 |
| | 1,2,3,7,8,9-HexaCDF | 1000 |
| | 2,3,4,6,7,8-HexaCDF | 1000 |
| | 1,2,3,4,6,7,8-HeptaCDD | 2000 |
| | 1,2,3,4,6,7,8-HeptaCDF | 2000 |
| | 1,2,3,4,7,8,9-HeptaCDF | 2000 |
| | OctaCDD | 2000 |
| | OctaCDF | 2000 |

JIS Methods K0311 and K0312 Dioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------------------|---|-----------------------------|
| NEW EDF-5327-H | Modified JIS Dioxin/Furan Calibration Solutions [CS1H-CS6H] | Set of 6 x 0.2 mL in nonane |
| NEW EDF-5327-H-E | Modified JIS Dioxin/Furan Calibration Solutions [CS0.2H-CS7H] | Set of 8 x 0.2 mL in nonane |
| NEW EDF-5327-CS0.2H | Modified JIS Dioxin/Furan Calibration Solution [CS0.2H] | 0.2 mL in nonane |
| NEW EDF-5327-CS1H | Modified JIS Dioxin/Furan Calibration Solution [CS1H] | 0.2 mL in nonane |
| NEW EDF-5327-CS2H | Modified JIS Dioxin/Furan Calibration Solution [CS2H] | 0.2 mL in nonane |
| NEW EDF-5327-CS3H | Modified JIS Dioxin/Furan Calibration Solution [CS3H] | 0.2 mL in nonane |
| NEW EDF-5327-CS4H | Modified JIS Dioxin/Furan Calibration Solution [CS4H] | 0.2 mL in nonane |
| NEW EDF-5327-CS5H | Modified JIS Dioxin/Furan Calibration Solution [CS5H] | 0.2 mL in nonane |
| NEW EDF-5327-CS6H | Modified JIS Dioxin/Furan Calibration Solution [CS6H] | 0.2 mL in nonane |
| NEW EDF-5327-CS7H | Modified JIS Dioxin/Furan Calibration Solution [CS7H] | 0.2 mL in nonane |

| All concentrations are in ng/mL (ppb) | | | | | | | | |
|---|--------|------|------|------|------|------|------|------|
| Unlabeled | CS0.2H | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H | CS7H |
| 2,3,7,8-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 2,3,7,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,3,6,8-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,3,6,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,3,7,9-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,8,9-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,7,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,8,9-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-PentaCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,3,7,8-PentaCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 2,3,4,7,8-PentaCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 |
| 1,2,3,4,7,8-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,4,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,6,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,7,8,9-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 2,3,4,6,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 |
| OctaCDD | 0.1 | 0.5 | 2.5 | 10 | 50 | 250 | 1000 | 2500 |
| OctaCDF | 0.1 | 0.5 | 2.5 | 10 | 50 | 250 | 1000 | 2500 |
| Labeled | | | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| OctaCDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |

JIS Methods K0311 and K0312 Dioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|--|---|------------------|
| NEW EDF-5328 | Modified JIS Dioxin/Furan Cleanup Spike | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 50 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | |
| OctaCDF (¹³ C ₁₂ , 99%) | 100 | |
| NEW EDF-5329 Modified JIS Dioxin/Furan Syringe Spike 1.2 mL in nonane | | |
| Labeled (ng/mL) | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7-PentaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | |

JIS Methods K0311 and K0312 Dioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------------------|---|-----------------------------|
| NEW EDF-5331-L-E | Modified JIS Low Concentration Dioxin/Furan Calibration Solutions [CS0.2L-CS5L] | Set of 6 x 0.2 mL in nonane |
| NEW EDF-5331-CS0.2L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS0.2L] | 0.2 mL in nonane |
| NEW EDF-5331-CS1L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS1L] | 0.2 mL in nonane |
| NEW EDF-5331-CS2L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS2L] | 0.2 mL in nonane |
| NEW EDF-5331-CS3L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS3L] | 0.2 mL in nonane |
| NEW EDF-5331-CS4L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS4L] | 0.2 mL in nonane |
| NEW EDF-5331-CS5L | Modified JIS Low Concentration Dioxin/Furan Calibration Solution [CS5L] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | CS0.2L | CS1L | CS2L | CS3L | CS4L | CS5L |
|---|--------|------|------|------|------|------|
| 2,3,7,8-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,3,6,8-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,3,7,9-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,8,9-TetraCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,3,7,8-PentaCDD | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,3,4,7,8-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,6,7,8-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,7,8,9-HexaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| OctaCDD | 0.1 | 0.5 | 2.5 | 10 | 50 | 250 |
| 2,3,7,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,3,6,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,7,8-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,8,9-TetraCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,3,7,8-PentaCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 2,3,4,7,8-PentaCDF | 0.02 | 0.1 | 0.5 | 2.0 | 10 | 50 |
| 1,2,3,4,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,6,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 2,3,4,6,7,8-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,7,8,9-HexaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.04 | 0.2 | 1.0 | 4.0 | 20 | 100 |
| OctaCDF | 0.1 | 0.5 | 2.5 | 10 | 50 | 250 |
| Labeled | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,7-PentaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| OctaCDD (¹³ C ₁₂ , 99%) | 4.0 | 4.0 | 4.0 | 4.0 | 4.0 | 4.0 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4-TetraCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 | 2.0 |
| OctaCDF (¹³ C ₁₂ , 99%) | 4.0 | 4.0 | 4.0 | 4.0 | 4.0 | 4.0 |

JIS Methods K0311 and K0312 Dioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| NEW EDF-5332 | Modified JIS Low Concentration Cleanup Spike | 1.2 mL in nonane |
| Labeled | | (ng/mL) |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 20 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 20 |
| NEW EDF-5333 | Modified JIS Low Concentration Syringe Spike | 1.2 mL in nonane |
| Labeled | | (ng/mL) |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,7-PentaCDD (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 |

JIS Methods K0311 and K0312 Dioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------|--|-----------------------------|
| EDF-5187 | JIS Dioxin/Furan Calibration Solutions [ST1-ST5] | Set of 5 x 0.2 mL in nonane |
| EDF-5187-ST1 | JIS Dioxin/Furan Calibration Solution [ST1] | 0.2 mL in nonane |
| EDF-5187-ST2 | JIS Dioxin/Furan Calibration Solution [ST2] | 0.2 mL in nonane |
| EDF-5187-ST3 | JIS Dioxin/Furan Calibration Solution [ST3] | 0.2 mL in nonane |
| EDF-5187-ST4 | JIS Dioxin/Furan Calibration Solution [ST4] | 0.2 mL in nonane |
| EDF-5187-ST5 | JIS Dioxin/Furan Calibration Solution [ST5] | 0.2 mL in nonane |

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | |
|---|-----|-----|-----|-----|-----|
| Unlabeled | ST1 | ST2 | ST3 | ST4 | ST5 |
| 2,3,7,8-TetraCDD | 0.2 | 1 | 5 | 20 | 100 |
| 1,2,3,7,8-PentaCDD | 0.2 | 1 | 5 | 20 | 100 |
| 1,2,3,4,7,8-HexaCDD | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,6,7,8-HexaCDD | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,7,8,9-HexaCDD | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.4 | 2 | 10 | 40 | 200 |
| OctaCDD | 1 | 5 | 25 | 100 | 500 |
| 2,3,7,8-TetraCDF | 0.2 | 1 | 5 | 20 | 100 |
| 1,2,3,7,8-PentaCDF | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,4,7,8-PentaCDF | 0.2 | 1 | 5 | 20 | 100 |
| 1,2,3,4,7,8-HexaCDF | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,6,7,8-HexaCDF | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,7,8,9-HexaCDF | 0.4 | 2 | 10 | 40 | 200 |
| 2,3,4,6,7,8-HexaCDF | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.4 | 2 | 10 | 40 | 200 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.4 | 2 | 10 | 40 | 200 |
| OctaCDF | 1 | 5 | 25 | 100 | 500 |
| Labeled | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 |
| OctaCDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 |

US EPA Method 8280 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EDF-2519-A | Method 8280 Calibration Solutions [CC1-CC5] | Set of 5 x 0.2 mL in nonane |
| EDF-2519-1 | Method 8280 Calibration Solution [CC1] | 0.2 mL in nonane |
| EDF-2519-2 | Method 8280 Calibration Solution [CC2] | 0.2 mL in nonane |
| EDF-2519-3 | Method 8280 Calibration and Verification Solution [CC3] | 0.2 mL in nonane |
| EDF-2519-4 | Method 8280 Calibration Solution [CC4] | 0.2 mL in nonane |
| EDF-2519-5 | Method 8280 Calibration Solution [CC5] | 0.2 mL in nonane |

| Unlabeled | All concentrations are in ng/µL (ppm) | | | | |
|---|---------------------------------------|-------|------|-----|------|
| | CC1 | CC2 | CC3 | CC4 | CC5 |
| 2,3,7,8-TetraCDD | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 2,3,7,8-TetraCDF | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-PentaCDF | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-PentaCDD | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 2,3,4,7,8-PentaCDF | — | — | 0.5 | — | — |
| 1,2,3,4,7,8-HexaCDF | — | — | 1.25 | — | — |
| 1,2,3,6,7,8-HexaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,7,8-HexaCDD | — | — | 1.25 | — | — |
| 1,2,3,6,7,8-HexaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,7,8,9-HexaCDD | — | — | 1.25 | — | — |
| 2,3,4,6,7,8-HexaCDF | — | — | 1.25 | — | — |
| 1,2,3,7,8,9-HexaCDF | — | — | 1.25 | — | — |
| 1,2,3,4,7,8,9-HeptaCDF | — | — | 1.25 | — | — |
| 1,2,3,4,6,7,8-HeptaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| OctaCDD | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| OctaCDF | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| Labeled | | | | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| OctaCDD (¹³ C ₁₂ , 99%) | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | — | — | 0.25 | — | — |

NEW EDF-8280-KIT Method 8280 "Starter Kit" * 1 Kit

Contains one each of the following items:

| | |
|------------|---|
| EDF-2519-A | Method 8280 Calibration Solutions [CS1-CS5] |
| EDF-2520 | Method 8280 Internal Standard Solution |
| ED-2521 | Method 8280 Recovery Standard Solution |
| ED-2522 | Method 8280 Cleanup Standard Solution |
| EDF-2523 | Method 8280 Matrix Spiking Solution |

*Modified Method 8280 "Starter Kit" also available

US EPA Method 8280 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EDF-4095 | Modified Method 8280 Calibration Solutions [CS1-CS5] (All 17 toxic congeners at all five levels) | Set of 5 × 0.2 mL in nonane |
| EDF-4095-1 | Modified Method 8280 Calibration Standard [CS1] | 0.2 mL in nonane |
| EDF-4095-2 | Modified Method 8280 Calibration Standard [CS2] | 0.2 mL in nonane |
| EDF-4095-3 | Modified Method 8280 Calibration Standard [CS3] | 0.2 mL in nonane |
| EDF-4095-4 | Modified Method 8280 Calibration Standard [CS4] | 0.2 mL in nonane |
| EDF-4095-5 | Modified Method 8280 Calibration Standard [CS5] | 0.2 mL in nonane |

All concentrations are in ng/µL (ppm)

| Unlabeled | CS1 | CS2 | CS3 | CS4 | CS5 |
|---|------|-------|------|------|------|
| 2,3,7,8-TetraCDD | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 2,3,7,8-TetraCDF | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-PentaCDD | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,7,8-PentaCDF | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 2,3,4,7,8-PentaCDD | 0.1 | 0.25 | 0.5 | 1.0 | 2.0 |
| 1,2,3,4,7,8-HexaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,7,8-HexaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,6,7,8-HexaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,6,7,8-HexaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,7,8,9-HexaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,7,8,9-HexaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 2,3,4,6,7,8-HexaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.25 | 0.625 | 1.25 | 2.5 | 5.0 |
| OctaCDD | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| OctaCDF | 0.5 | 1.25 | 2.5 | 5.0 | 10.0 |
| Labeled | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 0.25 | 0.25 | 0.25 | 0.25 | 0.25 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 0.5 | 0.5 | 0.5 | 0.5 | 0.5 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1.0 | 1.0 | 1.0 | 1.0 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | 1.0 | 1.0 | 1.0 | 1.0 | 1.0 |

| | | |
|----------|--|------------------|
| EDF-2681 | Supplemental Internal Standard Solution (not required by US EPA Method 8280) | 1.2 mL in nonane |
|----------|--|------------------|

| Labeled | (ng/µL) |
|---|---------|
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 5 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 5 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 5 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 10 |
| OctaCDF (¹³ C ₁₂ , 99%) | 10 |

US EPA Method 8280 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|------------------|
| EDF-2520 | Method 8280 Internal Standard Solution | 1.2 mL in nonane |
| | Labeled | (ng/µL) |
| | 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 5 |
| | 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 5 |
| | 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 5 |
| | 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 10 |
| | OctaCDD (¹³ C ₁₂ , 99%) | 10 |
| ED-2521 | Method 8280 Recovery Standard Solution | 1.2 mL in nonane |
| | Labeled | (ng/µL) |
| | 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 5 |
| | 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 5 |
| ED-2522 | Method 8280 Cleanup Standard Solution | 1.2 mL in nonane |
| | Labeled | (ng/µL) |
| | 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 5 |
| EDF-2523 | Method 8280 Matrix Spiking Solution | 1.2 mL in nonane |
| | Unlabeled | (ng/µL) |
| | 2,3,7,8-TetraCDD | 2.5 |
| | 2,3,7,8-TetraCDF | 2.5 |
| | 1,2,3,7,8-PentaCDF | 6.25 |
| | 1,2,3,7,8-PentaCDD | 6.25 |
| | 1,2,3,6,7,8-HexaCDF | 6.25 |
| | 1,2,3,6,7,8-HexaCDD | 6.25 |
| | 1,2,3,4,6,7,8-HeptaCDF | 6.25 |
| | 1,2,3,4,6,7,8-HeptaCDD | 6.25 |
| | OctaCDD | 12.5 |
| | OctaCDF | 12.5 |
| EDF-4096 | Modified Method 8280 Matrix Spiking Solution (all 17 toxic congeners) | 1.2 mL in nonane |
| | Unlabeled | (ng/µL) |
| | 2,3,7,8-TetraCDD | 2.5 |
| | 2,3,7,8-TetraCDF | 2.5 |
| | 1,2,3,7,8-PentaCDF | 6.25 |
| | 1,2,3,7,8-PentaCDD | 6.25 |
| | 2,3,4,7,8-PentaCDF | 6.25 |
| | 1,2,3,4,7,8-HexaCDD | 6.25 |
| | 1,2,3,4,7,8-HexaCDF | 6.25 |
| | 1,2,3,6,7,8-HexaCDD | 6.25 |
| | 1,2,3,6,7,8-HexaCDF | 6.25 |
| | 1,2,3,7,8,9-HexaCDD | 6.25 |
| | 1,2,3,7,8,9-HexaCDF | 6.25 |
| | 2,3,4,6,7,8-HexaCDF | 6.25 |
| | 1,2,3,4,6,7,8-HeptaCDD | 6.25 |
| | 1,2,3,4,6,7,8-HeptaCDF | 6.25 |
| | 1,2,3,4,7,8,9-HeptaCDF | 6.25 |
| | OctaCDD | 12.5 |
| | OctaCDF | 12.5 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|---------------|---|-----------------------------|
| EDF-5443 | Two-Column Dioxin and Furan and PCB Calibration Solutions [CS1H-CS6H] | Set of 6 x 0.2 mL in nonane |
| EDF-5443-CS1H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS1H] | 0.2 mL in nonane |
| EDF-5443-CS2H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS2H] | 0.2 mL in nonane |
| EDF-5443-CS3H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS3H] | 0.2 mL in nonane |
| EDF-5443-CS4H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS4H] | 0.2 mL in nonane |
| EDF-5443-CS5H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS5H] | 0.2 mL in nonane |
| EDF-5443-CS6H | Two-Column Dioxin and Furan and PCB Calibration Solution [CS6H] | 0.2 mL in nonane |

| Unlabeled | IUPAC | <i>All concentrations are in ng/mL (ppb)</i> | | | | | |
|--------------------------|-------|--|------|------|------|------|------|
| | | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
| 2,3,7,8-TetraCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,3,6,8-TetraCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,7,8-TetraCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,8,9-TetraCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,3,7,8-PentaCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,4,7,8-PentaCDF | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,3,4,7,8-HexaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 2,3,4,6,7,8-HexaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,4,7,8,9-HeptaCDF | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| OctaCDF | | 0.5 | 2.5 | 10 | 50 | 250 | 1000 |
| 2,3,7,8-TetraCDD | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,3,6,8-TetraCDD | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,3,7,9-TetraCDD | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,8,9-TetraCDD | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,3,7,8-PentaCDD | | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 |
| 1,2,3,4,7,8-HexaCDD | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDD | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDD | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDD | | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 |
| OctaCDD | | 0.5 | 2.5 | 10 | 50 | 250 | 1000 |
| 3,4,4',5-TetraCB | 81 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 3,3',4,4'-TetraCB | 77 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 3,3',4,4',5-PentaCB | 126 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2',3,4,4',5-PentaCB | 123 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,4,4',5-PentaCB | 114 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.2 | 0.5 | 2.0 | 10 | 50 | 200 |

(continued on next page)

Dioxin and Furan Plus PCB Standard Mixtures

(continued from previous page)

| Labeled | IUPAC | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
|---|-------|------|------|------|------|------|------|
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 | 10 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-----------------------|---|------------------|
| EDF-5444 | Two-Column Dioxin and Furan and PCB Cleanup Spike | 0.6 mL in nonane |
| NEW EDF-5444-C | Two-Column Dioxin and Furan and PCB Cleanup Spike with 1,3,6,8-TetraCDD | 0.6 mL in nonane |

| Labeled | IUPAC | EDF-5444 (ng/mL) | EDF-5444-C (ng/mL) |
|---|-------|---------------------|-----------------------|
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | | — | 1000 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 2000 | 2000 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 1000 | 1000 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 2000 | 2000 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 | 1000 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 | 1000 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 | 1000 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 | 1000 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 | 1000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 | 1000 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 | 1000 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 | 1000 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 | 1000 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 | 1000 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | 1000 |

| | | |
|--------------|---|------------------|
| EDF-5431 | Two-Column Dioxin and Furan Syringe Spike | 1.2 mL in nonane |
| EDF-5431-20X | Two-Column Dioxin and Furan Syringe Spike | 1.2 mL in nonane |

| Labeled | EDF-5431 (ng/mL) | EDF-5431-20X (ng/mL) |
|---|---------------------|-------------------------|
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | 1000 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) | 50 | 1000 |
| 1,2,3,4,6,7-HexaCDD (¹³ C ₁₂ , 99%) | 50 | 1000 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | 1000 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------|--------------------------------|------------------|
| EC-5325 | Modified JIS PCB Syringe Spike | 1.2 mL in nonane |
| EC-5325-0.2X | Modified JIS PCB Syringe Spike | 10 mL in nonane |
| EC-5325-20X | Modified JIS PCB Syringe Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5325 (ng/mL) | EC-5325-0.2X (ng/mL) | EC-5325-20X (ng/mL) |
|---|-------|--------------------|-------------------------|------------------------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 50 | 10 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 50 | 10 | 1000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 50 | 10 | 1000 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 50 | 10 | 1000 |

| | | |
|-------------|---------------------------------|------------------|
| EC-5326 | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |
| EC-5326-20X | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5326 (ng/mL) | EC-5326-20X (ng/mL) |
|--|-------|--------------------|------------------------|
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 50 | 1000 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|--|------------------------------|
| NEW EDF-5524 | Dioxin/Furan/CP-PCB Calibration Series [CDC1-CDC7] | Set of 7 x 0.25 mL in nonane |
| NEW EDF-5524-0.1 | Dioxin/Furan/CP-PCB Calibration Series [CDC0.1] | 0.25 mL in nonane |
| NEW EDF-5524-0.3 | Dioxin/Furan/CP-PCB Calibration Series [CDC0.3] | 0.25 mL in nonane |
| NEW EDF-5524-1 | Dioxin/Furan/CP-PCB Calibration Series [CDC1] | 0.25 mL in nonane |
| NEW EDF-5524-2 | Dioxin/Furan/CP-PCB Calibration Series [CDC2] | 0.25 mL in nonane |
| NEW EDF-5524-3 | Dioxin/Furan/CP-PCB Calibration Series [CDC3] | 0.25 mL in nonane |
| NEW EDF-5524-4 | Dioxin/Furan/CP-PCB Calibration Series [CDC4] | 0.25 mL in nonane |
| NEW EDF-5524-5 | Dioxin/Furan/CP-PCB Calibration Series [CDC5] | 0.25 mL in nonane |
| NEW EDF-5524-6 | Dioxin/Furan/CP-PCB Calibration Series [CDC6] | 0.25 mL in nonane |
| NEW EDF-5524-7 | Dioxin/Furan/CP-PCB Calibration Series [CDC7] | 0.25 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CDC0.1 | CDC0.3 | CDC1 | CDC2 | CDC3 | CDC4 | CDC5 | CDC6 | CDC7 |
|---|-------|--------|--------|------|------|------|------|------|------|------|
| 2,3,7,8-TetraCDD | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 2,3,7,8-TetraCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,7,8-PentaCDD | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,7,8-PentaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 2,3,4,7,8-PentaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,4,7,8-HexaCDD | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,4,7,8-HexaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,6,7,8-HexaCDD | | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 1,2,3,6,7,8-HexaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,7,8,9-HexaCDD | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,7,8,9-HexaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 2,3,4,6,7,8-HexaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD | | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF | | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 1,2,3,4,7,8,9-HeptaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| OctaCDD | | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 | 300 | 1000 |
| OctaCDF | | 0.001 | 0.003 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 |
| 3,3',4,4'-TetraCB | 77 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 3,4,4',5-TetraCB | 81 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 3,3',4,4',5-PentaCB | 126 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.01 | 0.03 | 0.1 | 0.3 | 1 | 3 | 10 | 30 | 100 |
| Labeled | | | | | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 3,3',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 80 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 | 30 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------------|--|---------------------------|
| NEW EDF-5525 | Dioxin/Furan/CP-PCB Internal Standard | 5.8 mL in methanol/nonane |
| NEW EDF-5525-100X-1.2 | Dioxin/Furan/CP-PCB Internal Standard 100X Stock | 1.2 mL in methanol/nonane |

| Labeled | IUPAC | EDF-5525 (pg/µL) | EDF-5525-100X-1.2 (pg/µL) |
|---|-------|---------------------|------------------------------|
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 1.5 | 150 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 1.5 | 150 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 1.5 | 150 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 5 | 500 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1.5 | 150 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1.5 | 150 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1.5 | 150 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1.5 | 150 |

| | | |
|------------------------------|--|---|
| NEW EDF-5526 | Dioxin/Furan/CP-PCB Recovery Standard | 5.8 mL in 88% hexane/ 10% nonane/2% dodecane |
| NEW EDF-5526-100X-1.2 | Dioxin/Furan/CP-PCB Recovery Standard 100X Stock | 1.2 mL in 88% hexane/ 10% nonane/2% dodecane |

| Labeled | IUPAC | EDF-5526 (pg/µL) | EDF-5526-100X-1.2 (pg/µL) |
|---|-------|---------------------|------------------------------|
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 0.5 | 50 |
| 3,3',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 80 | 1.5 | 150 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 0.5 | 50 |
| 2,4,6,8-TetraBDF (¹³ C ₁₂ , 99%) | | 1.0 | 100 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|----------------------------------|----------------------------------|
| NEW EDF-5538 | Dioxin/Furan/CP-PCB PAR Standard | 10 mL in nonane/toluene/isoctane |
| <hr/> | | |
| Unlabeled | IUPAC | (pg/ μ L) |
| 2,3,7,8-TetraCDD | | 20 |
| 2,3,7,8-TetraCDF | | 20 |
| 1,2,3,7,8-PentaCDD | | 20 |
| 1,2,3,7,8-PentaCDF | | 20 |
| 2,3,4,7,8-PentaCDF | | 20 |
| 1,2,3,4,7,8-HexaCDD | | 20 |
| 1,2,3,4,7,8-HexaCDF | | 20 |
| 1,2,3,6,7,8-HexaCDD | | 200 |
| 1,2,3,6,7,8-HexaCDF | | 20 |
| 1,2,3,7,8,9-HexaCDD | | 20 |
| 1,2,3,7,8,9-HexaCDF | | 20 |
| 2,3,4,6,7,8-HexaCDD | | 200 |
| 1,2,3,4,6,7,8-HeptaCDD | | 200 |
| 1,2,3,4,6,7,8-HeptaCDF | | 200 |
| 1,2,3,4,7,8,9-HeptaCDF | | 20 |
| OctaCDD | | 2000 |
| OctaCDF | | 20 |
| 3,3',4,4'-TetraCB | 77 | 200 |
| 3,4,4',5-TetraCB | 81 | 200 |
| 3,3',4,4',5-PentaCB | 126 | 200 |
| 3,3',4,4',5,5'-HexaCB | 169 | 200 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount | |
|---|---|--|-----------------------------|
| EDF-4144 | Internal Standard for Dioxin, Furan and PCB in Tissue | 750 µL in nonane | |
| Labeled | IUPAC | (ng/mL) | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 25.0 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 25.0 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 25.0 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 25.0 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 25.0 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 60.0 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 60.0 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 60.0 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 60.0 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| OctaCDD (¹³ C ₁₂ , 99%) | | 125 | |
| OctaCDF (¹³ C ₁₂ , 99%) | | 125 | |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 24.0 | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 24.0 | |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 36.0 | |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 48.0 | |
| EDF-4145 | Recovery Standard for Dioxin, Furan and PCB in Tissue | 750 µL in nonane | |
| Labeled | IUPAC | (ng/mL) | |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 25.0 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 62.5 | |
| 3,3',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 80 | 48.0 | |
| ES-5321 | Multi-Analyte Recovery Spiking Standard | 10 mL in 88% hexane/ 2% dodecane/10% nonane | |
| NEW ES-5321-200X-1.2 | Multi-Analyte Recovery Spiking Standard | 1.2 mL in nonane | |
| Labeled | IUPAC | ES-5321 (ng/mL) | ES-5321-200X-1.2 (ng/mL) |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 2.5 | 500 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 10.0 | 2000 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 7.5 | 1500 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 | 7.5 | 1500 |

Dioxin and Furan Plus PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|----------------------|------------------|
| EDF-5393 | Dioxin Cleanup Spike | 1.2 mL in nonane |

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 20 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 40 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 20 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 20 | |
| OctaCDF (¹³ C ₁₂ , 99%) | 40 | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 20 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 20 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 20 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 20 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 20 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 20 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 20 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 20 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 20 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 20 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 20 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 20 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 20 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 20 |

| | | |
|-------------------------|----------------------|------------------|
| EF-5394 | Dioxin Syringe Spike | 1.2 mL in nonane |
| NEW EF-5394-0.5X | Dioxin Syringe Spike | 1.2 mL in nonane |

| Labeled | EDF-5394 (ng/μL) | EDF-5394-0.5X (ng/μL) |
|---|---------------------|--------------------------|
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 20 | 10 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | 20 | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 20 | 10 |

| | | |
|----------|-----------------------|------------------|
| EDF-5395 | Dioxin Sampling Spike | 1.2 mL in nonane |
|----------|-----------------------|------------------|

| Labeled | IUPAC | (ng/μL) |
|--|-------|---------|
| 1,2,3,4-TetraCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 50 |

Non-2,3,7,8-Containing Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount | | | | |
|-------------|--|-----------------------------|--|--|--|--|
| EDF-5392 | Dioxin/Furan Calibration Solutions [CS1-CS6] | Set of 6 × 0.2 mL in nonane | | | | |
| EDF-5392-1 | Dioxin/Furan Calibration Solution [CS1] | 0.2 mL in nonane | | | | |
| EDF-5392-2 | Dioxin/Furan Calibration Solution [CS2] | 0.2 mL in nonane | | | | |
| EDF-5392-3 | Dioxin/Furan Calibration Solution [CS3] | 0.2 mL in nonane | | | | |
| EDF-5392-4 | Dioxin/Furan Calibration Solution [CS4] | 0.2 mL in nonane | | | | |
| EDF-5392-5 | Dioxin/Furan Calibration Solution [CS5] | 0.2 mL in nonane | | | | |
| EDF-5392-6 | Dioxin/Furan Calibration Solution [CS6] | 0.2 mL in nonane | | | | |

| Unlabeled | | All concentrations are in ng/mL (ppb) | | | | | |
|---|--|---------------------------------------|-----|-----|-----|-----|------|
| | | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
| 2,3,7,8-TetraCDD | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,3,6,8-TetraCDD | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,3,7,9-TetraCDD | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,8,9-TetraCDD | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,3,7,8-PentaCDD | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,3,4,7,8-HexaCDD | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDD | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDD | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDD | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| OctaCDD | | 0.5 | 2.5 | 10 | 50 | 250 | 1000 |
| 2,3,7,8-TetraCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,3,6,8-TetraCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,7,8-TetraCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,8,9-TetraCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,3,7,8-PentaCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,4,7,8-PentaCDF | | 0.1 | 0.5 | 2 | 10 | 50 | 200 |
| 1,2,3,4,7,8-HexaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 2,3,4,6,7,8-HexaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| 1,2,3,4,7,8,9-HeptaCDF | | 0.2 | 1 | 4 | 20 | 100 | 400 |
| OctaCDF | | 0.5 | 2.5 | 10 | 50 | 250 | 1000 |
| Cleanup | | | | | | | |
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDF (¹³ C ₁₂ , 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Syringe | | | | | | | |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) (alternate) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4-TetraCDF (¹³ C ₁₂ , 99%) | | 10 | 10 | 10 | 10 | 10 | 10 |

Dioxin and Furan Food/Feed/QQQ Standard Mixtures

| Catalog No. | Compound | Amount |
|---------------------------|---|------------------------------|
| NEW EDF-5554 | Dioxin and Furan Food / Feed / QQQ Calibration Series [CS1-CS8] * | Set of 8 x 0.25 mL in nonane |
| NEW EDF-5554-L | Dioxin and Furan Food / Feed / QQQ Calibration Series [CS1-CS7] * | Set of 7 x 0.25 mL in nonane |
| NEW EDF-5554-H | Dioxin and Furan Food / Feed / QQQ Calibration Series [CS2-CS8] * | Set of 7 x 0.25 mL in nonane |
| NEW EDF-5554-CS0.4 | Dioxin and Furan Food / Feed / QQQ Calibration Series [CS0.4] * | 0.25 mL in nonane |
| NEW EDF-5554-CS0.2 | Dioxin and Furan Food / Feed / QQQ Calibration Series [CS0.2] * | 0.25 mL in nonane |

Individual calibration solutions and user-defined combinations are available. Please inquire.

All concentrations are in pg/ μ L (ppb)

| Unlabeled | CS0.2 | CS0.4 | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 |
|--|-------|-------|------|-----|-----|-----|-----|-----|-----|------|
| 2,3,7,8-TetraCDD | 0.01 | 0.02 | 0.05 | 0.1 | 0.5 | 2 | 5 | 25 | 50 | 200 |
| 2,3,7,8-TetraCDF | 0.01 | 0.02 | 0.05 | 0.1 | 0.5 | 2 | 5 | 25 | 50 | 200 |
| 1,2,3,7,8-PentaCDD | 0.01 | 0.02 | 0.05 | 0.1 | 0.5 | 2 | 5 | 25 | 50 | 200 |
| 1,2,3,7,8-PentaCDF | 0.01 | 0.02 | 0.05 | 0.1 | 0.5 | 2 | 5 | 25 | 50 | 200 |
| 2,3,4,7,8-PentaCDD | 0.01 | 0.02 | 0.05 | 0.1 | 0.5 | 2 | 5 | 25 | 50 | 200 |
| 1,2,3,4,7,8-HexaCDD | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDD | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDD | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,4,7,8-HexaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,6,7,8-HexaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,7,8,9-HexaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 2,3,4,6,7,8-HexaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.02 | 0.04 | 0.1 | 0.2 | 1 | 4 | 10 | 50 | 100 | 400 |
| OctaCDD | 0.05 | 0.1 | 0.25 | 0.5 | 2.5 | 10 | 25 | 125 | 250 | 1000 |
| OctaCDF | 0.05 | 0.1 | 0.25 | 0.5 | 2.5 | 10 | 25 | 125 | 250 | 1000 |
| Extraction | | | | | | | | | | |
| 2,3,7,8-TetraCDD ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDF ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,6,7,8-HexaCDD ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,7,8,9-HexaCDD ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,7,8-HexaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,6,7,8-HexaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,7,8,9-HexaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,7,8-HeptaCDD ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,7,8-HeptaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,7,8-HeptaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,7,8,9-HeptaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| OctaCDD ($^{13}\text{C}_{12}$, 99%) | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| OctaCDF ($^{13}\text{C}_{12}$, 99%) | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 | 50 |
| Cleanup | | | | | | | | | | |
| 1,2,3,4-TetraCDD ($^{13}\text{C}_6$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Injection | | | | | | | | | | |
| 1,2,3,4-TetraCDD ($^{13}\text{C}_{12}$, 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,9-HexaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4,6,8,9-HeptaCDF ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |

*In production; please inquire for final specifications.

Dioxin and Furan Food/Feed/QQQ Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|---|------------------|
| NEW EDF-5555 | Dioxin and Furan Food / Feed / QQQ Extraction Standard Solution * | 1.2 mL in nonane |
| NEW EDF-5555-10X | Dioxin and Furan Food / Feed / QQQ Extraction Standard Solution * | 1.2 mL in nonane |

| Labeled | EDF-5555 (pg/µL) | EDF-5555-10X (pg/µL) |
|---|---------------------|-------------------------|
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 100 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 20 | 200 |
| OctaCDD (¹³ C ₁₂ , 99%) | 50 | 500 |
| OctaCDF (¹³ C ₁₂ , 99%) | 50 | 500 |

| | | |
|-------------------------|--|------------------|
| NEW EDF-5556 | Dioxin and Furan Food/Feed/QQQ Injection Standard Solution * | 1.2 mL in nonane |
| NEW EDF-5556-10X | Dioxin and Furan Food/Feed/QQQ Injection Standard Solution * | 1.2 mL in nonane |

| Labeled | EDF-5556 (pg/µL) | EDF-5556-10X (pg/µL) |
|---|---------------------|-------------------------|
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 20 | 200 |
| 1,2,3,4,6,9-HexaCDF (¹³ C ₁₂ , 99%) | 40 | 400 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 40 | 400 |

| | | |
|-------------------------|--|------------------|
| NEW EDF-5574 | Dioxin and Furan Food/Feed/QQQ Cleanup Standard Solution * | 1.2 mL in nonane |
| NEW EDF-5574-10X | Dioxin and Furan Food/Feed/QQQ Cleanup Standard Solution * | 1.2 mL in nonane |

| Labeled | EDF-5574 (pg/µL) | EDF-5574-10X (pg/µL) |
|--|---------------------|-------------------------|
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 20 | 200 |

| | | |
|---------------------|--|------------------|
| NEW EDF-5557 | Dioxin and Furan Food/Feed/QQQ PAR Standard Solution * | 1.2 mL in nonane |
|---------------------|--|------------------|

| Unlabeled | (pg/µL) | Unlabeled | (pg/µL) |
|---------------------|---------|------------------------|---------|
| 2,3,7,8-TetraCDD | 500 | 1,2,3,4,6,7,8-HeptaCDD | 1000 |
| 2,3,7,8-TetraCDF | 500 | 1,2,3,4,6,7,8-HeptaCDF | 1000 |
| 1,2,3,7,8-PentaCDD | 500 | 1,2,3,4,7,8,9-HeptaCDF | 1000 |
| 1,2,3,7,8-PentaCDF | 500 | OctaCDD | 2500 |
| 2,3,4,7,8-PentaCDF | 500 | OctaCDF | 2500 |
| 1,2,3,4,7,8-HexaCDD | 1000 | | |
| 1,2,3,6,7,8-HexaCDD | 1000 | | |
| 1,2,3,7,8,9-HexaCDD | 1000 | | |
| 1,2,3,4,7,8-HexaCDF | 1000 | | |
| 1,2,3,6,7,8-HexaCDF | 1000 | | |
| 1,2,3,7,8,9-HexaCDF | 1000 | | |
| 2,3,4,6,7,8-HexaCDF | 1000 | | |

*In production; please inquire for final specifications.

Two-Column Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|---------------|---|-----------------------------|
| EDF-5429-6H | Two-Column Dioxin and Furan Calibration Solutions [CS1H-CS6H] | Set of 6 x 0.2 mL in nonane |
| EDF-5429-7H | Two-Column Dioxin and Furan Calibration Solutions [CS1H-CS7H] | Set of 7 x 0.2 mL in nonane |
| EDF-5429-CS1H | Two-Column Dioxin and Furan Calibration Solution [CS1H] | 0.2 mL in nonane |
| EDF-5429-CS2H | Two-Column Dioxin and Furan Calibration Solution [CS2H] | 0.2 mL in nonane |
| EDF-5429-CS3H | Two-Column Dioxin and Furan Calibration Solution [CS3H] | 0.2 mL in nonane |
| EDF-5429-CS4H | Two-Column Dioxin and Furan Calibration Solution [CS4H] | 0.2 mL in nonane |
| EDF-5429-CS5H | Two-Column Dioxin and Furan Calibration Solution [CS5H] | 0.2 mL in nonane |
| EDF-5429-CS6H | Two-Column Dioxin and Furan Calibration Solution [CS6H] | 0.2 mL in nonane |
| EDF-5429-CS7H | Two-Column Dioxin and Furan Calibration Solution [CS7H] | 0.2 mL in nonane |
| EDF-5429-CS8H | Two-Column Dioxin and Furan Calibration Solution [CS8H] | 0.2 mL in nonane |

| All concentrations are in ng/mL (ppb) | | | | | | | | |
|---|------|------|------|------|------|------|------|------|
| Unlabeled | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H | CS7H | CS8H |
| 2,3,7,8-TetraCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,3,6,8-TetraCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,7,8-TetraCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,8,9-TetraCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,3,7,8-PentaCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 2,3,4,7,8-PentaCDF | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,3,4,7,8-HexaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,6,7,8-HexaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 2,3,4,6,7,8-HexaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,7,8,9-HexaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,4,6,7,8-HeptaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,4,7,8,9-HeptaCDF | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| OctaCDF | 0.5 | 2.5 | 10 | 50 | 250 | 1000 | 2500 | 5000 |
| 2,3,7,8-TetraCDD | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,3,6,8-TetraCDD | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,3,7,9-TetraCDD | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,8,9-TetraCDD | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,3,7,8-PentaCDD | 0.1 | 0.5 | 2.0 | 10 | 50 | 200 | 500 | 1000 |
| 1,2,3,4,7,8-HexaCDD | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,6,7,8-HexaCDD | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,7,8,9-HexaCDD | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| 1,2,3,4,6,7,8-HeptaCDD | 0.2 | 1.0 | 4.0 | 20 | 100 | 400 | 1000 | 2000 |
| OctaCDD | 0.5 | 2.5 | 10 | 50 | 250 | 1000 | 2500 | 5000 |
| Labeled | | | | | | | | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |

Two-Column Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|---|---|--|
| EDF-5430 | Two-Column Dioxin and Furan Cleanup Spike | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 50 | |
| OctaCDF (¹³ C ₁₂ , 99%) | 100 | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 50 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 100 | |
| | | |
| EDF-5431 | Two-Column Dioxin and Furan Syringe Spike | 1.2 mL in nonane |
| EDF-5431-20X | Two-Column Dioxin and Furan Syringe Spike | 1.2 mL in nonane |
| | | |
| Labeled | | |
| | | EDF-5431 (ng/mL) |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 50 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) | | 50 |
| 1,2,3,4,6,7-HexaCDD (¹³ C ₁₂ , 99%) | | 50 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 50 |
| | | EDF-5431-20X (ng/mL) |
| 1,2,7,8-TetraCDF (¹³ C ₁₂ , 99%) | | 1000 |
| 1,2,3,4,6-PentaCDF (¹³ C ₁₂ , 99%) | | 1000 |
| 1,2,3,4,6,7-HexaCDD (¹³ C ₁₂ , 99%) | | 1000 |
| 1,2,3,4,6,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | | 1000 |
| | | |
| NEW | ED-5432 | Two-Column Dioxin and Furan Sampling Spike |
| | | |
| Labeled | | (ng/mL) |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | | 50 |

Isotope-Labeled Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|---|------------------|
| EDF-4067 | Tetra-OctaCDD and CDF Standard Solution (2,3,7,8-isomers) | 1.2 mL in nonane |
| EDF-4067-A | Tetra-OctaCDD and CDF Standard Solution (2,3,7,8-isomers excluding 1,2,3,7,8,9-HexaCDD) | 1.2 mL in nonane |
| EDF-5041 | Non-2,3,7,8-Containing PCDF Cleanup Standard | 1.2 mL in nonane |
| NEW EDF-5041-20 | Non-2,3,7,8-Containing PCDF Cleanup Standard 1/20 Dilution | 5 mL in nonane |

| Labeled | EDF-4067 (ng/mL) | EDF-4067-A (ng/mL) | EDF-5041 (ng/mL) | EDF-5041-20 (ng/mL) |
|---|---------------------|-----------------------|---------------------|------------------------|
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | — | 2000 | 100 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | 1000 | 2000 | 100 |
| OctaCDD (¹³ C ₁₂ , 99%) | 1000 | 2000 | 4000 | 200 |
| OctaCDF (¹³ C ₁₂ , 99%) | 1000 | 2000 | 4000 | 200 |

| | | |
|--------------|---|------------------|
| EDF-5304 | Dioxin and Furan Cleanup Spike | 1.2 mL in nonane |
| EDF-5174-40X | 1,3,6,8-TCDD/F Containing Cleanup Spike | 1.2 mL in nonane |

| Labeled | EDF-5304 (ng/mL) | EDF-5174-40X (ng/mL) |
|---|---------------------|-------------------------|
| 1,3,6,8-TetraCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,7,8,9-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 1000 | 200 |
| OctaCDD (¹³ C ₁₂ , 99%) | 2000 | 400 |
| 1,3,6,8-TetraCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,7,8,9-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| 1,2,3,4,7,8,9-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | 200 |
| OctaCDF (¹³ C ₁₂ , 99%) | 2000 | 400 |

Isotope-Labeled Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|-----------------------------|
| EDF-4964-A | JIS Dioxin/Furan Type 1 Cleanup Standard Solution | 0.5 mL in nonane |
| Labeled (ng/mL) | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 2,3,4,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,4,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2000 | |
| 2,3,4,6,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 2000 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 2000 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 4000 | |
| OctaCDF (¹³ C ₁₂ , 99%) | 4000 | |
| | | |
| EDF-957 | Carbon-13 Quantifying Cocktail (2,3,7,8-PCDD/PCDF isomers) | Set of 3 × 0.4 mL in nonane |
| Labeled (ng/mL) | | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,7,8-PentaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,7,8-PentaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,6,7,8-HexaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,7,8-HexaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,6,7,8-HeptaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| 1,2,3,4,6,7,8-HeptaCDF (¹³ C ₁₂ , 99%) | 1000 | |
| OctaCDD (¹³ C ₁₂ , 99%) | 1000 | |
| OctaCDF (¹³ C ₁₂ , 99%) | 1000 | |

Unlabeled Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------------------|--|--------------------|
| NEW EDF-5493 | Dioxin/Furan Native Mix All 17 2,3,7,8 Isomers + 1,3,6,8-TetraCDD/F + 1,3,7,9-TetraCDD | 1.2 mL in nonane |
| Unlabeled (ng/mL) | | |
| 1,3,6,8-TetraCDD | 1000 | |
| 2,3,7,8-TetraCDD | 1000 | |
| 1,3,7,9-TetraCDD | 1000 | |
| 1,2,3,7,8-PentaCDD | 1000 | |
| 1,2,3,4,7,8-HexaCDD | 2000 | |
| 1,2,3,6,7,8-HexaCDD | 2000 | |
| 1,2,3,7,8,9-HexaCDD | 2000 | |
| 1,2,3,4,6,7,8-HeptaCDD | 2000 | |
| OctaCDD | 5000 | |
| 1,3,6,8-TetraCDF | 1000 | |
| 2,3,7,8-TetraCDF | 1000 | |
| 1,2,3,7,8-PentaCDF | 1000 | |
| 2,3,4,7,8-PentaCDF | 2000 | |
| 1,2,3,4,7,8-HexaCDF | 2000 | |
| 1,2,3,6,7,8-HexaCDF | 2000 | |
| 1,2,3,7,8,9-HexaCDF | 2000 | |
| 2,3,4,6,7,8-HexaCDF | 2000 | |
| 1,2,3,4,7,8,9-HeptaCDF | 2000 | |
| OctaCDF | 5000 | |
| | | |
| ED-906B-5 | TCDD-HpCDD Standard Solution (B) (2,3,7,8 isomers) | 1.2 mL in nonane |
| ED-906B-25 | TCDD-HpCDD Standard Solution (B) (2,3,7,8 isomers) | 1.2 mL in nonane |
| | | |
| Unlabeled | | |
| | ED-906B-5 (ng/mL) | ED-906B-25 (ng/mL) |
| 2,3,7,8-TetraCDD | 5000 | 25,000 |
| 1,2,3,7,8-PentaCDD | 5000 | 25,000 |
| 1,2,3,4,7,8-HexaCDD | 5000 | 25,000 |
| 1,2,3,6,7,8-HexaCDD | 5000 | 25,000 |
| 1,2,3,7,8,9-HexaCDD | 5000 | 25,000 |
| 1,2,3,4,6,7,8-HeptaCDD | 5000 | 25,000 |
| | | |
| EF-909B-5 | TCDF-HpCDF Standard Solution (B) (2,3,7,8 isomers) | 1.2 mL in nonane |
| EF-909B-25 | TCDF-HpCDF Standard Solution (B) (2,3,7,8 isomers) | 1.2 mL in nonane |
| | | |
| Unlabeled | | |
| | ED-909B-5 (ng/mL) | ED-909B-25 (ng/mL) |
| 2,3,7,8-TetraCDF | 5000 | 25,000 |
| 1,2,3,7,8-PentaCDF | 5000 | 25,000 |
| 2,3,4,7,8-PentaCDF | 5000 | 25,000 |
| 1,2,3,4,7,8-HexaCDF | 5000 | 25,000 |
| 1,2,3,6,7,8-HexaCDF | 5000 | 25,000 |
| 1,2,3,7,8,9-HexaCDF | 5000 | 25,000 |
| 2,3,4,6,7,8-HexaCDF | 5000 | 25,000 |
| 1,2,3,4,6,7,8-HeptaCDF | 5000 | 25,000 |
| 1,2,3,4,7,8,9-HeptaCDF | 5000 | 25,000 |

Unlabeled Dioxin and Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|-------------------------------|----------------|
| ED-4135 | Chlorodioxin Mix – High | 1 mL in nonane |
| Unlabeled | | (ng/mL) |
| 2-MonoCDD | | 5000 |
| 2,8-DiCDD | | 5000 |
| 2,3,7-TriCDD | | 5000 |
| 2,3,7,8-TetraCDD | | 5000 |
| 1,2,3,7,8-PentaCDD | | 5000 |
| 1,2,3,4,7,8-HexaCDD | | 5000 |
| 1,2,3,6,7,8-HexaCDD | | 5000 |
| 1,2,3,7,8,9-HexaCDD | | 5000 |
| 1,2,3,4,6,7,8-HeptaCDD | | 5000 |
| OctaCDD | | 5000 |
| EF-4134 | Chlorodibenzofuran Mix – High | 1 mL in nonane |
| Unlabeled | | (ng/mL) |
| 2-MonoCDF | | 5000 |
| 2,4-DiCDF | | 5000 |
| 2,4,6-TriCDF | | 5000 |
| 2,3,7,8-TetraCDF | | 5000 |
| 1,2,3,7,8-PentaCDF | | 5000 |
| 2,3,4,7,8-PentaCDF | | 5000 |
| 1,2,3,4,7,8-HexaCDF | | 5000 |
| 1,2,3,6,7,8-HexaCDF | | 5000 |
| 2,3,4,6,7,8-HexaCDF | | 5000 |
| 1,2,3,7,8,9-HexaCDF | | 5000 |
| 1,2,3,4,6,7,8-HeptaCDF | | 5000 |
| 1,2,3,4,7,8,9-HeptaCDF | | 5000 |
| OctaCDF | | 5000 |

Window Defining Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| EDF-4147 | PCDD/PentaCDF Window Defining and Isomer Specificity Mix (DB-5 and DB-225 Columns) | 1.2 mL in nonane |
| Unlabeled (ng/mL) | | |
| 1,3,6,8-TetraCDD | 200 | |
| 1,2,8,9-TetraCDD | 200 | |
| 2,3,7,8-TetraCDD | 200 | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 200 | |
| 1,2,3,7/1,2,3,8-TetraCDD | 200 | |
| 1,2,3,9-TetraCDD | 200 | |
| 1,3,6,8-TetraCDF | 200 | |
| 1,2,8,9-TetraCDF | 200 | |
| 2,3,7,8-TetraCDF | 200 | |
| 2,3,7,8-TetraCDF (¹³ C ₁₂ , 99%) | 200 | |
| 2,3,4,7-TetraCDF | 200 | |
| 1,2,3,9-TetraCDF | 200 | |
| 1,2,4,6,8/1,2,4,7,9-PentaCDD | 200 | |
| 1,2,3,8,9-PentaCDD | 200 | |
| 1,3,4,6,8-PentaCDF | 200 | |
| 1,2,3,8,9-PentaCDF | 200 | |
| 1,2,4,6,7,9/1,2,4,6,8,9-HexaCDD | 200 | |
| 1,2,3,4,6,7-HexaCDD | 200 | |
| 1,2,3,4,6,8-HexaCDF | 200 | |
| 1,2,3,4,8,9-HexaCDF | 200 | |
| 1,2,3,4,6,7,9-HeptaCDD | 200 | |
| 1,2,3,4,6,7,8-HeptaCDF | 200 | |
| 1,2,3,4,6,7,8-HeptaCDD | 200 | |
| 1,2,3,4,7,8,9-HeptaCDF | 200 | |
| | | |
| ED-1732-S | TCDD-HpCDD Window Defining Mixture (DB-5) | 0.5 mL in nonane |
| Unlabeled (ng/mL) | | |
| 1,3,6,8-TetraCDD | 800 | |
| 1,2,8,9-TetraCDD | 800 | |
| 1,2,4,6,8/1,2,4,7,9-PentaCDD | 800 | |
| 1,2,3,8,9-PentaCDD | 800 | |
| 1,2,3,4,6,7-HexaCDD | 800 | |
| 1,2,4,6,7,9/1,2,4,6,8,9-HexaCDD | 800 | |
| 1,2,3,4,6,7,8-HeptaCDD | 800 | |
| 1,2,3,4,6,7,9-HeptaCDD | 800 | |
| | | |
| EF-1731-S | TCDF-HpCDF Window Defining Mixture (DB-5) | 0.5 mL in nonane |
| Unlabeled (ng/mL) | | |
| 1,3,6,8-TetraCDF | 800 | |
| 1,2,8,9-TetraCDF | 800 | |
| 1,3,4,6,8-PentaCDF | 800 | |
| 1,2,3,8,9-PentaCDF | 800 | |
| 1,2,3,4,6,8-HexaCDF | 800 | |
| 1,2,3,4,8,9-HexaCDF | 800 | |
| 1,2,3,4,6,7,8-HeptaCDF | 800 | |
| 1,2,3,4,7,8,9-HeptaCDF | 800 | |

Column Performance Mixtures

| Catalog No. | Compound | Amount |
|---|---|------------------|
| ED-908 | TCDD Column Performance Solution Mixture | 1.2 mL in nonane |
| Unlabeled (ng/mL) | | |
| 1,2,3,4-TetraCDD | 10 | |
| 1,2,3,7/1,2,3,8-TetraCDD | 10 | |
| 1,2,7,8-TetraCDD | 10 | |
| 1,4,7,8-TetraCDD | 10 | |
| 2,3,7,8-TetraCDD | 10 | |
| | | |
| ED-935-A | Modified TCDD Column Performance Check Solution | 0.5 mL in nonane |
| Unlabeled (ng/mL) | | |
| 2,3,7,8-TetraCDD | 100 | |
| 1,2,3,4-TetraCDD | 100 | |
| 1,4,7,8-TetraCDD | 100 | |
| 1,2,3,7/1,2,3,8-TetraCDD | 100 | |
| 1,2,7,8-TetraCDD | 200 | |
| 2,3,7,8-TetraCDD (¹³ C ₁₂ , 99%) | 250 | |
| 2,3,7,8-TetraCDD (³⁷ Cl ₄ , 96%) | 7 | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 500 | |

Instrument Performance Mixture

| Catalog No. | Compound | Amount |
|---|---|------------------|
| ED-5461 | 1,2,3,4-TCDD Instrument Performance Check | 0.1 mL in nonane |
| | | |
| Unlabeled (pg/mL) | | |
| 1,2,3,4-TetraCDD | 100 | |
| Labeled | | |
| 1,2,3,4-TetraCDD (¹³ C ₁₂ , 99%) | 100 | |

Bromodioxin/Furan Calibration Solutions

| Catalog No. | Compound | Amount |
|-------------|--|-----------------------------|
| EDF-5407 | Bromodioxin/Furan Calibration Standard Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EDF-5407-1 | Bromodioxin/Furan Calibration Standard Solution [CS1] | 0.2 mL in nonane |
| EDF-5407-2 | Bromodioxin/Furan Calibration Standard Solution [CS2] | 0.2 mL in nonane |
| EDF-5407-3 | Bromodioxin/Furan Calibration Standard Solution [CS3] | 0.2 mL in nonane |
| EDF-5407-4 | Bromodioxin/Furan Calibration Standard Solution [CS4] | 0.2 mL in nonane |
| EDF-5407-5 | Bromodioxin/Furan Calibration Standard Solution [CS5] | 0.2 mL in nonane |

| Unlabeled | <i>All concentrations are in ng/mL (ppb)</i> | | | | |
|------------------------|---|-----|------|-----|-----|
| | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,3,7,8-TetraBDD | 0.1 | 0.4 | 2.0 | 10 | 50 |
| 1,2,3,7,8-PentaBDD | 0.2 | 0.8 | 4.0 | 20 | 100 |
| 1,2,3,4,7,8-HexabDD | 0.6 | 2.4 | 12.0 | 60 | 300 |
| 1,2,3,6,7,8-HexabDD | 0.6 | 2.4 | 12.0 | 60 | 300 |
| 1,2,3,7,8,9-HexabDD | 0.6 | 2.4 | 12.0 | 60 | 300 |
| 1,2,3,4,6,7,8-HeptaBDD | 0.75 | 3.0 | 15.0 | 75 | 375 |
| OctaBDD | 1.0 | 4.0 | 20.0 | 100 | 500 |
| 2,3,7,8-TetraBDF | 0.2 | 0.8 | 4.0 | 20 | 100 |
| 2,4,6,8-TetraBDF | 0.2 | 0.8 | 4.0 | 20 | 100 |
| 1,2,3,7,8-PentaBDF | 0.4 | 1.6 | 8.0 | 40 | 200 |
| 2,3,4,7,8-PentaBDF | 0.4 | 1.6 | 8.0 | 40 | 200 |
| 1,2,3,4,7,8-HexaBDF | 0.6 | 2.4 | 12.0 | 60 | 300 |
| 1,2,3,4,6,7,8-HeptaBDF | 0.75 | 3.0 | 15.0 | 75 | 375 |
| OctaBDF | 1.0 | 4.0 | 20.0 | 100 | 500 |
| Labeled | CS1 | CS2 | CS3 | CS4 | CS5 |
| | 2,3,7,8-TetraBDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 1,2,3,7,8-PentaBDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 1,2,3,4,7,8-HexabDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 |
| | 1,2,3,6,7,8-HexabDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 |
| | 1,2,3,7,8,9-HexabDD (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 |
| | 1,2,3,4,6,7,8-HeptaBDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 |
| | OctaBDD (¹³ C ₁₂ , 99%) | 150 | 150 | 150 | 150 |
| | 2,3,7,8-TetraBDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 2,4,6,8-TetraBDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 1,2,3,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 2,3,4,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 |
| | 1,2,3,4,7,8-HexaBDF (¹³ C ₁₂ , 99%) | 50 | 50 | 50 | 50 |
| | 1,2,3,4,6,7,8-HeptaBDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 |
| | OctaBDF (¹³ C ₁₂ , 99%) | 150 | 150 | 150 | 150 |

Bromodioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|---------------------------------------|
| EDF-5408 | Bromodioxin/Furan Cleanup Spike | 0.5 mL in nonane |
| <hr/> | | |
| | Labeled | (ng/mL) |
| | 2,3,7,8-TetraBDD (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,7,8-PentaBDD (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 250 |
| | 1,2,3,6,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 250 |
| | 1,2,3,4,6,7,8-HeptaBDD (¹³ C ₁₂ , 99%) | 500 |
| | OctaBDD (¹³ C ₁₂ , 99%) | 750 |
| | 2,3,7,8-TetraBDF (¹³ C ₁₂ , 99%) | 100 |
| | 2,3,4,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 100 |
| | 1,2,3,4,7,8-HexaBDF (¹³ C ₁₂ , 99%) | 250 |
| | 1,2,3,4,6,7,8-HeptaBDF (¹³ C ₁₂ , 99%) | 500 |
| | OctaBDF (¹³ C ₁₂ , 99%) | 750 |
| <hr/> | | |
| EDF-5409-A | Bromodioxin/Furan Syringe Spike | 1.2 mL in nonane:toluene |
| <hr/> | | |
| | Labeled | (ng/mL) |
| | 1,2,3,7,8,9-HexaBDD (¹³ C ₁₂ , 99%) | 500 |
| | 1,2,3,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 200 |
| <hr/> | | |
| EF-5410 | Bromodioxin/Furan Sampling Spike | 1.2 mL in nonane |
| <hr/> | | |
| | Labeled | (ng/mL) |
| | 2,4,6,8-TetraBDF (¹³ C ₁₂ , 99%) | 200 |
| <hr/> | | |
| NEW | EDF-5517 | Bromodioxin/Furan Native PAR Solution |
| <hr/> | | |
| | Unlabeled | (ng/mL) |
| | 2,3,7,8-TetraBDD | 50 |
| | 1,2,3,7,8-PentaBDD | 100 |
| | 1,2,3,4,7,8-HexaBDD | 300 |
| | 1,2,3,6,7,8-HexaBDD | 300 |
| | 1,2,3,7,8,9-HexaBDD | 300 |
| | 1,2,3,4,6,7,8-HeptaBDD | 375 |
| | OctaBDD | 500 |
| | 2,3,7,8-TetraBDF | 100 |
| | 2,4,6,8-TetraBDF | 100 |
| | 1,2,3,7,8-PentaBDF | 200 |
| | 2,3,4,7,8-PentaBDF | 200 |
| | 1,2,3,4,7,8-HexaBDF | 300 |
| | 1,2,3,4,6,7,8-HeptaBDF | 375 |
| | OctaBDF | 500 |

Bromodioxin/Furan Calibration Solutions

| Catalog No. | Compound | Amount |
|--------------|--|-----------------------------|
| EDF-5381 | PBDD/F Calibration Solutions [CS1-CS7] | Set of 7 x 0.2 mL in nonane |
| EDF-5381-CS1 | PBDD/F Calibration Solution [CS1] | 0.2 mL in nonane |
| EDF-5381-CS2 | PBDD/F Calibration Solution [CS2] | 0.2 mL in nonane |
| EDF-5381-CS3 | PBDD/F Calibration Solution [CS3] | 0.2 mL in nonane |
| EDF-5381-CS4 | PBDD/F Calibration Solution [CS4] | 0.2 mL in nonane |
| EDF-5381-CS5 | PBDD/F Calibration Solution [CS5] | 0.2 mL in nonane |
| EDF-5381-CS6 | PBDD/F Calibration Solution [CS6] | 0.2 mL in nonane |
| EDF-5381-CS7 | PBDD/F Calibration Solution [CS7] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 |
|---|------|-----|-----|-----|-----|-----|-----|
| 2,3,7,8-TetraBDD | 0.1 | 0.4 | 2 | 10 | 20 | 40 | — |
| 1,2,3,7,8-PentaBDD | 0.2 | 0.8 | 4 | 20 | 40 | 80 | — |
| 1,2,3,4,7,8-HexaBDD | 0.75 | 3 | 15 | 75 | 150 | 300 | — |
| 1,2,3,6,7,8-HexaBDD | 0.75 | 3 | 15 | 75 | 150 | 300 | — |
| 1,2,3,7,8,9-HexaBDD | 0.75 | 3 | 15 | 75 | 150 | 300 | — |
| OctaBDD | 1 | 4 | 20 | 100 | 200 | 400 | 800 |
| 2,3,7,8-TetraBDF | 0.5 | 2 | 10 | 50 | 100 | 200 | — |
| 2,4,6,8-TetraBDF | 0.5 | 2 | 10 | 50 | 100 | 200 | — |
| 1,2,3,7,8-PentaBDF | 0.5 | 2 | 10 | 50 | 100 | 200 | — |
| 2,3,4,7,8-PentaBDF | 0.5 | 2 | 10 | 50 | 100 | 200 | — |
| 1,2,3,4,7,8-HexaBDF | 0.75 | 3 | 15 | 75 | 150 | 300 | — |
| 1,2,3,4,6,7,8-HeptaBDF | 0.75 | 3 | 15 | 75 | 150 | 300 | 600 |
| OctaBDF | 1 | 4 | 20 | 100 | 200 | 400 | 800 |
| Labeled | | | | | | | |
| 2,3,7,8-TetraBDD (¹³ C ₁₂ , 99%) | 10 | 10 | 10 | 10 | 10 | 10 | — |
| 1,2,3,7,8-PentaBDD (¹³ C ₁₂ , 99%) | 20 | 20 | 20 | 20 | 20 | 20 | — |
| 1,2,3,4,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 75 | 75 | 75 | 75 | 75 | 75 | — |
| 1,2,3,6,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 75 | 75 | 75 | 75 | 75 | 75 | — |
| 1,2,3,7,8,9-HexaBDD (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | — |
| OctaBDD (¹³ C ₁₂ , 99%) | 225 | 225 | 225 | 225 | 225 | 225 | 225 |
| 2,3,7,8-TetraBDF (¹³ C ₁₂ , 99%) | 40 | 40 | 40 | 40 | 40 | 40 | — |
| 2,4,6,8-TetraBDF (¹³ C ₁₂ , 99%) | 40 | 40 | 40 | 40 | 40 | 40 | — |
| 1,2,3,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 40 | 40 | 40 | 40 | 40 | 40 | — |
| 2,3,4,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 40 | 40 | 40 | 40 | 40 | 40 | — |
| 1,2,3,4,7,8-HexaBDF (¹³ C ₁₂ , 99%) | 40 | 40 | 40 | 40 | 40 | 40 | — |
| 1,2,3,4,6,7,8-HeptaBDF (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | — |
| OctaBDF (¹³ C ₁₂ , 99%) | 225 | 225 | 225 | 225 | 225 | 225 | 225 |

Bromodioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|---|----------------------------|------------------------|
| EDF-5382 | PBDD/F Cleanup Spike | 0.5 mL in nonane |
| Labeled | | |
| | (ng/mL) | |
| 2,3,7,8-TetraBDD (¹³ C ₁₂ , 99%) | 50 | |
| 1,2,3,7,8-PentaBDD (¹³ C ₁₂ , 99%) | 100 | |
| 1,2,3,4,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 375 | |
| 1,2,3,6,7,8-HexaBDD (¹³ C ₁₂ , 99%) | 375 | |
| OctaBDD (¹³ C ₁₂ , 99%) | 1125 | |
| 2,3,7,8-TetraBDF (¹³ C ₁₂ , 99%) | 200 | |
| 2,3,4,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 200 | |
| 1,2,3,4,7,8-HexaBDF (¹³ C ₁₂ , 99%) | 375 | |
| 1,2,3,4,6,7,8-HeptaBDF (¹³ C ₁₂ , 99%) | 500 | |
| OctaBDF (¹³ C ₁₂ , 99%) | 1125 | |
| EDF-5383 | | |
| EDF-5383 | PBDD/F Syringe Spike Stock | 1.2 mL in nonane |
| EDF-5383-4X | PBDD/F Syringe Spike Stock | 1.2 mL in nonane |
| Labeled | | |
| | EDF-5383 (ng/mL) | EDF-5383-4X (ng/mL) |
| 1,2,3,7,8,9-HexaBDD (¹³ C ₁₂ , 99%) | 500 | 2000 |
| 1,2,3,7,8-PentaBDF (¹³ C ₁₂ , 99%) | 200 | 800 |
| EF-5384 | | |
| EF-5384 | PBDD/F Sampling Stock | 1.2 mL in nonane |
| EF-5384-4X | PBDD/F Sampling Stock | 1.2 mL in nonane |
| Labeled | | |
| | EF-5384 (ng/mL) | EF-5384-4X (ng/mL) |
| 2,4,6,8-TetraBDF (¹³ C ₁₂ , 99%) | 200 | 800 |

Bromodioxin/Furan Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------|--|------------------|
| ED-5073 | Brominated Dioxin/Furan Internal Standard | 1.2 mL in nonane |
| Labeled | | |
| | (ng/mL) | |
| | 2,3,7,8-TetraBDD (¹³ C ₁₂ , 99%) | 200 |
| | 1,2,3,6,7,8-HexabDD (¹³ C ₁₂ , 99%) | 50 |
| | 1,2,3,7,8,9-HexabDD (¹³ C ₁₂ , 99%) | 150 |
| | | |
| EDF-5059 | Polybrominated Dioxin and Furan Mixture | 1.2 mL in nonane |
| Unlabeled | | |
| | (ng/mL) | |
| | 2,3,7,8-TetraBDD | 1000 |
| | 1,2,3,7,8-PentaBDD | 1000 |
| | 1,2,3,4,7,8-HexabDD | 1000 |
| | 1,2,3,6,7,8-HexabDD | 1000 |
| | 1,2,3,7,8,9-HexabDD | 1000 |
| | OctabDD | 1000 |
| | 2,3,7,8-TetraBDF | 1000 |
| | 1,2,3,7,8-PentaBDF | 1000 |
| | 2,3,4,7,8-PentaBDF | 1000 |
| | 1,2,3,4,7,8-HexabDF | 1000 |
| | 1,2,3,4,6,7,8-HeptaBDF | 1000 |



From their first commercial use in the 1920s, it is estimated that over 1,500,000 tons (1.5×10^{21} ng) of PCBs were produced worldwide. PCBs were banned from production in most countries in the 1970s, though large-scale contamination of the environment remains a problem today. In 2001, PCBs were included in the original Stockholm Convention on Persistent Organic Pollutants.

PCB Nomenclature

In recent years, interest in PCB testing has centered around two different but related groups of congeners, one group based on congeners most commonly found in technical PCB products and thus in the environment, and another based on toxicity. To help researchers more easily find the products they need, PCB standards and standard mixtures have been categorized by these two groups: "WHO Dioxin-Like PCBs" – compounds that demonstrate dioxin-like activity and have been given toxic equivalence factors (TEFs) by the World Health Organization (WHO); and "WHO Non-Dioxin-Like (Marker/Indicator) PCBs" – compounds that are found in the highest concentrations in technical products and the environment, but have not been given TEFs by the WHO. Please note that CIL also offers a set of standard mixtures developed for the combined analysis of both groups. Look to CIL for even more developments in this area as we continue to work toward meeting the complete needs of researchers in environmental, food, water, and exposure analysis.

Unlabeled "CS" PCB Standards

CIL's primary business is isotopically labeled standards, but it is important to remember that the accuracy and precision of a quantitative analysis is dependent upon the accuracy and precision of the unlabeled (native) standards. In the past, CIL utilized commercially available native standards from multiple vendors for the confirmation of its isotopically labeled standards. It was discovered, however, that there can be substantial variability among the commercial native standards. Thus, CIL initiated the "CS" (certified) PCB standards program.

CIL prepares native certified standards using good laboratory practice (GLP). Individual, native crystalline PCB isomers (98%+ purity) are weighed in triplicate on a microbalance calibrated with NIST-traceable Class S weights and formulated to specific concentration. Triplicate analyses of each of the three solutions in isoctane are carried out using GC/MS. In order to establish statistical control, the relative standard deviation (RSD) of each solution must be less than 5%, and the RSD for the entire set of analyses for all three standards must be <5%. When these parameters have been met, the solutions are combined and the resulting solution analyzed again in triplicate by three chemists. If the RSD of these analyses is also <5%, the final product is the certified PCB standard. These 100 µg/mL solutions are highly accurate native standards for quantitation of PCBs. These standards are used in all CIL calibration series and native standard mixtures, and are used to validate all isotope-labeled standards from CIL.

Isotope-Labeled PCB Standards

CIL offers more than 50 individual ¹³C-labeled PCB standards to meet the growing needs of researchers utilizing isotope dilution mass spectrometry (IDMS). All ¹³C-labeled PCB standards are quantified against CIL-certified unlabeled PCB standards for utmost precision and accuracy.

High-Purity PCB Standards

CIL responds to the needs of the analytical community by providing high-purity PCB standards. As new instrumentation and methodologies drive detection limits lower, the presence of even very low levels of impurities in the labeled standards of other PCB congeners or polychlorinated dibenzo-*p*-dioxins and dibenzofurans (PCDD/Fs) becomes a hindrance to a laboratory's minimum-detection capabilities.

CIL has developed aggressive cleanup procedures and adopted much tighter quality-control specifications for the eight mono-ortho-substituted dioxin-like PCBs (DL-PCBs). These new specifications include extremely low allowances for ¹³C-non-ortho DL-PCBs, native content, other PCB congeners and PCDD/Fs.

High-Purity PCB Standards

Chemical Identity: Unambiguous identity by GC-MS,
¹H-NMR, ¹³C-NMR, and MP determination

Isotopic Enrichment: 99% by GC-MS

Chemical Purity: >98% by GC-MS, GC-ECD, and ¹H-NMR

- Native Content: <0.1% by GC-MS SI
- ¹³C-non-ortho DL-PCBs: <0.05% by GC-ECD vs. cal-curve, or HRGC-MS
- 17 (2,3,7,8) containing PCDD/Fs: <0.05% for each compound by HRGC-MS

Concentration: 40 ± 2 µg/mL by comparison assay vs. native "certified standard"

Uncertainty: Conforming to Eurachem/CITAC Guide
"Quantifying Uncertainty in Analytical Measurement"

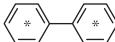
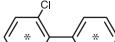
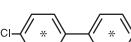
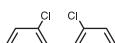
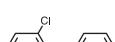
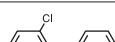
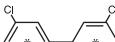
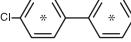
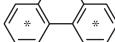
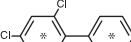
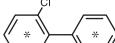
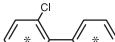
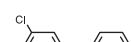
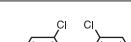
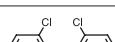
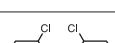
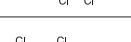
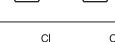
Comprehensive Native PCB Mixtures

CIL has produced a new calibration series and spiking solutions to go along with its two large mixes of unlabeled PCBs formulated entirely from its PCB certified standards individual stock solutions. The Comprehensive PCB Mixtures include all the WHO dioxin-like PCBs that have been assigned TEFs, the predominant congeners, and first- and last-eluting congeners from the mono-through-deca homologue groups. Since some pairs of these compounds coelute on certain columns, the Fully Resolved Native Mono-Deca PCB Mixture was formulated with no coeluting congeners under normal analytical conditions.

Mixed Bromo/Chlorobiphenyl Standards

CIL offers a selection of labeled and unlabeled mixed halogenated biphenyl standards and standard mixtures. While very limited research on these compounds has been done to date, they have been identified in environmental matrices and warrant further investigation.

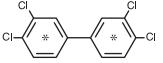
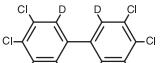
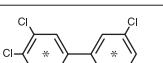
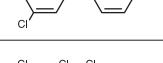
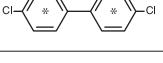
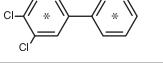
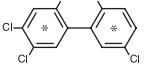
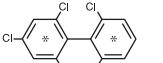
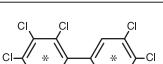
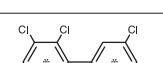
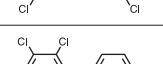
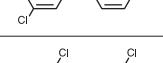
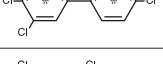
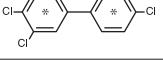
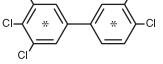
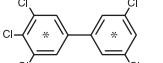
Isotope-Labeled Individual PCB Standards

| Catalog No. | Compound | IUPAC | Structure | Concentration | Amount |
|--|---|-------|--|------------------------|----------------|
| CLM-3235-1.2 | Biphenyl ($^{13}\text{C}_{12}$, 99%) | 0 |  | 100 µg/mL in nonane | 1.2 mL |
| EC-4908-3 EC-4908-1.2 | 2-Monochlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 1 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4990-3 EC-4990-1.2 | 4-Monochlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 3 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4911-3 EC-4911-1.2 | 2,2'-Dichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 4 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-5095-3 EC-5095-1.2 | 2,4'-Dichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 8 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4165-3 EC-4165-1.2 | 2,5-Dichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 9 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| NEW EC-5494-3 NEW EC-5494-1.2 | 3,3'-Dichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 11 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1402-3 EC-1402-1.2 | 4,4'-Dichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 15 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4909-3 EC-4909-1.2 | 2,2',6-Trichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 19 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| M EC-1413-3 EC-1413-1.2 | 2,4,4'-Trichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 28 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| NEW EC-5512-3 NEW EC-5512-1.2 | 2,4',5-Trichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 31 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4163-3 EC-4163-1.2 | 2,4',6-Trichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 32 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4901-3 EC-4901-1.2 | 3,4,4'-Trichlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 37 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1434-3 EC-1434-1.2 | 2,2',4,4'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 47 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| M EC-1424-3 EC-1424-1.2 | 2,2',5,5'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 52 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4912-3 EC-4912-1.2 | 2,2',6,6'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 54 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4078-3 EC-4078-1.2 | 2,3,4,4'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 60 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4914-3 EC-4914-1.2 | 2,3',4',5-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 70 |  | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |

M = marker PCB

DL = dioxin-like PCB

Isotope-Labeled Individual PCB Standards

| | Catalog No. | Compound | IUPAC | Structure | Concentration | Amount |
|-------------|--------------|---|-------|--|--|--------|
| DL | EC-1404-3 | 3,3',4,4'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 77 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1404-1.2 | | | | | 1.2 mL |
| DL | DLM-3063-3 | 3,3',4,4'-Tetrachlorobiphenyl (D_6 , 98%) | 77 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | DLM-3063-1.2 | | | | | 1.2 mL |
| M | EC-5048-3 | 3,3',4,5'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 79 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-5048-1.2 | | | | | 1.2 mL |
| DL | EC-1414-3 | 3,3',5,5'-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 80 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1414-1.2 | | | | | 1.2 mL |
| DL | EC-1412-3 | 3,4,4',5-Tetrachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 81 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1412-1.2 | | | | | 1.2 mL |
| M | EC-4929-3 | 2,2',3,4,4'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 85 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-4929-1.2 | | | | | 1.2 mL |
| M | EC-1428-3 | 2,2',3',4,5-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 97 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1428-1.2 | | | | | 1.2 mL |
| M | EC-1405-3 | 2,2',4,5,5'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 101 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1405-1.2 | | | | | 1.2 mL |
| DL | EC-4910-3 | 2,2',4,6,6'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 104 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-4910-1.2 | | | | | 1.2 mL |
| DL | EC-1420-3 | 2,3,3',4,4'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 105 |  | 40 ± 2 µg/mL in nonane *high purity | 3 mL |
| | EC-1420-1.2 | | | | | 1.2 mL |
| M | EC-1415-3 | 2,3,3',5,5'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 111 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1415-1.2 | | | | | 1.2 mL |
| DL | EC-4902-3 | 2,3,4,4',5-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 114 |  | 40 ± 2 µg/mL in nonane *high purity | 3 mL |
| | EC-4902-1.2 | | | | | 1.2 mL |
| M/DL | EC-1435-3 | 2,3',4,4',5-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 118 |  | 40 ± 2 µg/mL in nonane *high purity | 3 mL |
| | EC-1435-1.2 | | | | | 1.2 mL |
| DL | EC-4904-3 | 2',3,4,4',5-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 123 |  | 40 ± 2 µg/mL in nonane *high purity | 3 mL |
| | EC-4904-1.2 | | | | | 1.2 mL |
| DL | EC-1425-3 | 3,3',4,4',5-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 126 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1425-1.2 | | | | | 1.2 mL |
| M | EC-1421-3 | 3,3',4,5,5'-Pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 127 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1421-1.2 | | | | | 1.2 mL |
| M | EC-1411-3 | 2,2',3,3',4,4'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 128 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1411-1.2 | | | | | 1.2 mL |
| M | EC-1436-3 | 2,2',3,4,4',5-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 138 |  | 40 ± 2 µg/mL in nonane | 3 mL |
| | EC-1436-1.2 | | | | | 1.2 mL |

M = marker PCB**DL** = dioxin-like PCB

Isotope-Labeled Individual PCB Standards

| Catalog No. | Compound | IUPAC | Structure | Concentration | Amount |
|-----------------|--|-------|-----------|--|----------------|
| EC-1426-3 | 2,2',3,4,5,5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 141 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| M EC-1406-3 | 2,2',4,4',5,5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 153 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1406-1.2 | | | | | |
| EC-4167-3 | 2,2',4,4',6,6'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 155 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4167-1.2 | | | | | |
| DL EC-1422-3 | 2,3,3',4,4',5-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 156 | | 40 ± 2 µg/mL in nonane *high purity | 3 mL 1.2 mL |
| EC-1422-1.2 | | | | | |
| DL EC-4051-3 | 2,3,3',4,4',5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 157 | | 40 ± 2 µg/mL in nonane *high purity | 3 mL 1.2 mL |
| EC-4051-1.2 | | | | | |
| EC-5336-3 | 2,3,3',4,5,5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 159 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-5336-1.2 | | | | | |
| DL EC-4050-3 | 2,3',4,4',5,5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 167 | | 40 ± 2 µg/mL in nonane *high purity | 3 mL 1.2 mL |
| EC-4050-1.2 | | | | | |
| DL EC-1416-3 | 3,3',4,4',5,5'-Hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 169 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1416-1.2 | | | | | |
| EC-4905-3 | 2,2',3,3',4,4',5-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 170 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4905-1.2 | | | | | |
| EC-1417-3 | 2,2',3,3',5,5',6-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 178 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1417-1.2 | | | | | |
| M EC-1407-3 | 2,2',3,4,4',5,5'-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 180 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1407-1.2 | | | | | |
| NEW EC-5471-3 | 2,2',3,4,4',5,6'-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 182 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| NEW EC-5471-1.2 | | | | | |
| EC-4913-3 | 2,2',3,4',5,6,6'-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 188 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4913-1.2 | | | | | |
| DL EC-1409-3 | 2,3,3',4,4',5,5'-Heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 189 | | 40 ± 2 µg/mL in nonane *high purity | 3 mL 1.2 mL |
| EC-1409-1.2 | | | | | |
| EC-1418-3 | 2,2',3,3',4,4',5,5'-Octachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 194 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1418-1.2 | | | | | |
| EC-1408-3 | 2,2',3,3',5,5',6,6'-Octachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 202 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-1408-1.2 | | | | | |
| EC-4199-3 | 2,3,3',4,4',5,5',6-Octachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 205 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4199-1.2 | | | | | |
| EC-4900-3 | 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 206 | | 40 ± 2 µg/mL in nonane | 3 mL 1.2 mL |
| EC-4900-1.2 | | | | | |

M = marker PCB

DL = dioxin-like PCB

Isotope-Labeled Individual PCB Standards

| Catalog No. | Compound | IUPAC | Structure | Concentration | Amount |
|-------------|---|-------|-----------|------------------------|--------|
| EC-1419-3 | 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 208 | | 40 ± 2 µg/mL in nonane | 3 mL |
| EC-1419-1.2 | (¹³ C ₁₂ , 99%) | | | | 1.2 mL |
| EC-1410-3 | Decachlorobiphenyl (¹³ C ₁₂ , 99%) | 209 | | 40 ± 2 µg/mL in nonane | 3 mL |
| EC-1410-1.2 | | | | | 1.2 mL |
| EC-1410-10 | | | | | 10 mL |

Unlabeled Individual PCB "CS" Standards

| Catalog No. | Compound | Concentration | Amount |
|-------------|-----------------------|---------------------------------|----------------------------|
| PCB-1-CS | 2-Monochlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-3-CS | 4-Monochlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-4-CS | 2,2'-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-8-CS | 2,4'-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-9-CS | 2,5-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-10-CS | 2,6-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-11-CS | 3,3'-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-12-CS | 3,4-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-15-CS | 4,4'-Dichlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| NEW | PCB-17-CS | 2,2',4-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-18-CS | 2,2',5-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-19-CS | 2,2',6-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-28-CS | 2,4,4'-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-30-CS | 2,4,6-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-31-CS | 2,4',5-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-32-CS | 2,4',6-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-33-CS | 2',3,4-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-35-CS | 3,3',4-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-37-CS | 3,4,4'-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-38-CS | 3,4,5-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-39-CS | 3,4',5-Trichlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-44-CS | 2,2',3,5'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-47-CS | 2,2',4,4'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| NEW | PCB-49-CS | 2,2',4,5'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-52-CS | 2,2',5,5'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-54-CS | 2,2',6,6'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-57-CS | 2,3,3',5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-60-CS | 2,3,4,4'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-66-CS | 2,3',4,4'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-70-CS | 2,3',4',5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-74-CS | 2,4,4',5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-77-CS | 3,3',4,4'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-78-CS | 3,3',4,5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-79-CS | 3,3',4,5',5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-80-CS | 3,3',5,5'-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-81-CS | 3,4,4',5-Tetrachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-82-CS | 2,2',3,3',4-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| NEW | PCB-85-CS | 2,2',3,4,4'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-87-CS | 2,2',3,4,5'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-95-CS | 2,2',3,5',6-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-97-CS | 2,2',3',4,5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-99-CS | 2,2',4,4',5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-101-CS | 2,2',4,5,5'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-104-CS | 2,2',4,6,6'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-105-CS | 2,3,3',4,4'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |
| | PCB-110-CS | 2,3,3',4',6-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane |

Unlabeled Individual PCB "CS" Standards

| Catalog No. | Compound | Concentration | Amount |
|-------------|-----------------------------------|--|---------------------------|
| PCB-111-CS | 2,3,3',5,5'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isooctane | 1.2 mL |
| PCB-112-CS | 2,3,3',5,6-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-114-CS | 2,3,4,4',5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-118-CS | 2,3',4,4',5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-123-CS | 2',3,4,4',5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-126-CS | 3,3',4,4',5-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-127-CS | 3,3',4,5,5'-Pentachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| PCB-128-CS | 2,2',3,3',4,4'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane | 1.2 mL |
| NEW | PCB-132-CS | 2,2',3,3',4,6'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-138-CS | 2,2',3,4,4',5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-141-CS | 2,2',3,4,5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-146-CS | 2,2',3,4',5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-149-CS | 2,2',3,4',5',6-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-151-CS | 2,2',3,5,5',6-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-153-CS | 2,2',4,4',5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-155-CS | 2,2',4,4',6,6'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-156-CS | 2,3,3',4,4',5-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-157-CS | 2,3,3',4,4',5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-158-CS | 2,3,3',4,4',6-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-159-CS | 2,3,3',4,5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-162-CS | 2,3,3',4',5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-167-CS | 2,3',4,4',5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-169-CS | 3,3',4,4',5,5'-Hexachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-170-CS | 2,2',3,3',4,4',5-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| NEW | PCB-171-CS | 2,2',3,3',4,4',6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-172-CS | 2,2',3,3',4,5,5'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-174-CS | 2,2',3,3',4,5,6'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-177-CS | 2,2',3,3',4',5,6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-178-CS | 2,2',3,3',5,5',6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-180-CS | 2,2',3,4,4',5,5'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-182-CS | 2,2',3,4,4',5,6'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-183-CS | 2,2',3,4,4',5',6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-187-CS | 2,2',3,4',5,5',6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-188-CS | 2,2',3,4',5,6,6'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| NEW | PCB-189-CS | 2,3,3',4,4',5,5'-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-191-CS | 2,3,3',4,4',5',6-Heptachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-194-CS | 2,2',3,3',4,4',5,5'-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-195-CS | 2,2',3,3',4,4',5,6-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-196-CS | 2,2',3,3',4,4',5',6-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-198-CS | 2,2',3,3',4,5,5',6-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-199-CS | 2,2',3,3',4,5,6,6'-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-201-CS | 2,2',3,3',4,5,5',6'-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-202-CS | 2,2',3,3',5,5',6,6'-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-203-CS | 2,2',3,4,4',5,5',6-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| NEW | PCB-205-CS | 2,3,3',4,4',5,5',6-Octachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-206-CS | 2,2',3,3',4,4',5,5',6-Nonachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-208-CS | 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl | 100 ± 5 µg/mL in isoctane |
| | PCB-209-CS | Decachlorobiphenyl | 100 ± 5 µg/mL in isoctane |

Unlabeled PCB Standards

CIL also offers:

- All 209 PCBs in single weighed solutions at approximately 35 µg/mL in isoctane, with chemical purity >95%.
- All 209 PCBs in neat form; typically at nominal weights of 1 or 5 mg.

For details, please consult the CIL website at shop.isotope.com.

US EPA Method 1668A/B/C Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|-----------------------------|
| EC-4976 | Method 1668A/B/C Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EC-4976-0.2 | Method 1668A/B/C High Sensitivity Calibration Solution [CS0.2] (not included in EC-4976) | 0.2 mL in nonane |
| EC-4976-1 | Method 1668A/B/C Calibration Solution [CS1] | 0.2 mL in nonane |
| EC-4976-2 | Method 1668A/B/C Calibration Solution [CS2] | 0.2 mL in nonane |
| EC-4976-3 | Method 1668A/B/C Calibration Verification Solution [CS3] | 0.2 mL in nonane |
| EC-4976-3-4 | Method 1668A/B/C Calibration Verification Solution [CS3] | Set of 4 x 0.2 mL in nonane |
| EC-4976-4 | Method 1668A/B/C Calibration Solution [CS4] | 0.2 mL in nonane |
| EC-4976-5 | Method 1668A/B/C Calibration Solution [CS5] | 0.2 mL in nonane |

| All concentrations are in ng/mL (ppb) | | | | | | | |
|--|-------|-------|-----|-----|-----|-----|------|
| Native Toxics/LOC | IUPAC | CS0.2 | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2-MonoCB | 1 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 4-MonoCB | 3 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2'-DiCB | 4 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 4,4'-DiCB | 15 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',6-TriCB | 19 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,4,4'-TriCB | 37 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',6,6'-TetraCB | 54 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4'-TetraCB | 77 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,4,4',5-TetraCB | 81 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',4,6,6'-PentaCB | 104 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,4,4',5-PentaCB | 114 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2',3,4,4',5-PentaCB | 123 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4',5-PentaCB | 126 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',4,4',6,6'-HexaCB | 155 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,4',5,6,6'-HeptaCB | 188 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',5,5',6,6'-OctaCB | 202 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,3,3',4,4',5,5',6-OctaCB | 205 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| DecaCB | 209 | 0.2 | 1.0 | 5.0 | 50 | 400 | 2000 |
| Labeled Toxics/LOC/Window Defining | | | | | | | |
| 2-MonoCB (¹³ C ₁₂ , 99%) | 1 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2'-DiCB (¹³ C ₁₂ , 99%) | 4 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',6-TriCB (¹³ C ₁₂ , 99%) | 19 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,4,4'-TriCB (¹³ C ₁₂ , 99%) | 37 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',6,6'-TetraCB (¹³ C ₁₂ , 99%) | 54 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,6,6'-PentaCB (¹³ C ₁₂ , 99%) | 104 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',6,6'-HexaCB (¹³ C ₁₂ , 99%) | 155 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 100 | 100 | 100 | 100 | 100 | 100 |

(continued on next page)

US EPA Method 1668A/B/C Standard Mixtures

(continued from previous page)

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | | |
|---|-------|-------|-----|-----|-----|-----|-----|
| Labeled Toxics/LOC/Window Defining | IUPAC | CS0.2 | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,2',3,4',5,6,6'-HeptaCB (¹³ C ₁₂ , 99%) | 188 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',5,5',6,6'-OctaCB (¹³ C ₁₂ , 99%) | 202 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4',5,5',6-OctaCB (¹³ C ₁₂ , 99%) | 205 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 100 | 100 | 100 | 100 | 100 | 100 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 100 | 100 | 100 | 100 | 100 | 100 |
| Labeled Cleanup | | | | | | | |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 100 | 100 | 100 | 100 | 100 | 100 |
| Labeled Injection Internal | | | | | | | |
| 2,5-DiCB (¹³ C ₁₂ , 99%) | 9 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 100 | 100 | 100 | 100 | 100 | 100 |

NEW EC-1668A/B-KIT Method 1668A/B/C "Starter Kit"

1 Kit

Contains one each of the following items:

| | |
|---------|---|
| EC-4976 | Method 1668A/B/C Calibration Solutions [CS1-CS5] |
| EC-4977 | Method 1668A/B/C Labeled Toxics/LOC/Window Defining Solution |
| EC-4978 | Method 1668A/B/C Labeled Cleanup Standard Solution |
| EC-4979 | Method 1668A/B/C Labeled Injection Internal Standard Solution |
| EC-4989 | Method 1668A/B/C Native Toxics/LOC Solution |

US EPA Method 1668A/B/C Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|------------------|
| EC-4977 | Method 1668A/B/C Labeled Toxics/LOC/Window Defining Solution | 1.2 mL in nonane |
| EC-4977-5 | Method 1668A/B/C Labeled Toxics/LOC/Window Defining Solution | 5 mL in nonane |

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2-MonoCB (¹³ C ₁₂ , 99%) | 1 | 1000 |
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 1000 |
| 2,2'-DiCB (¹³ C ₁₂ , 99%) | 4 | 1000 |
| 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 1000 |
| 2,2',6-TriCB (¹³ C ₁₂ , 99%) | 19 | 1000 |
| 3,4,4'-TriCB (¹³ C ₁₂ , 99%) | 37 | 1000 |
| 2,2',6,6'-TetraCB (¹³ C ₁₂ , 99%) | 54 | 1000 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 |
| 2,2',4,6,6'-PentaCB (¹³ C ₁₂ , 99%) | 104 | 1000 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 155 | 1000 |
| 2,2',4,4',6,6'-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 |
| 2,2',3,4',5,6,6'-HeptaCB (¹³ C ₁₂ , 99%) | 188 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 |
| 2,2',3,3',5,5',6,6'-OctaCB (¹³ C ₁₂ , 99%) | 202 | 1000 |
| 2,3,3',4,4',5,5',6-OctaCB (¹³ C ₁₂ , 99%) | 205 | 1000 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 1000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 1000 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 1000 |

| EC-4978 | Method 1668A/B/C Labeled Cleanup Standard Solution | 1.2 mL in nonane |
|---|--|------------------|
| Labeled | IUPAC | (ng/mL) |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 1000 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 1000 |

| EC-4979 | Method 1668A/B/C Labeled Injection Internal Standard Solution | 1.2 mL in nonane |
|---|---|------------------|
| Labeled | IUPAC | (ng/mL) |
| 2,5-DiCB (¹³ C ₁₂ , 99%) | 9 | 5000 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 5000 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 5000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 5000 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 5000 |

US EPA Method 1668A/B/C Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------------|---|------------------|
| EC-4989 | Method 1668A/B/C Native Toxics/LOC Solution | 1.2 mL in nonane |
| Unlabeled | | |
| 2-MonoCB | IUPAC | (ng/mL) |
| 2-MonoCB | 1 | 2000 |
| 4-MonoCB | 3 | 2000 |
| 2,2'-DiCB | 4 | 2000 |
| 4,4'-DiCB | 15 | 2000 |
| 2,2',6-TriCB | 19 | 2000 |
| 3,4,4'-TriCB | 37 | 2000 |
| 2,2',6,6'-TetraCB | 54 | 2000 |
| 3,3',4,4'-TetraCB | 77 | 2000 |
| 3,4,4',5-TetraCB | 81 | 2000 |
| 2,2',4,6,6'-PentaCB | 104 | 2000 |
| 2,3,3',4,4'-PentaCB | 105 | 2000 |
| 2,3,4,4',5-PentaCB | 114 | 2000 |
| 2,3',4,4',5-PentaCB | 118 | 2000 |
| 2',3,4,4',5-PentaCB | 123 | 2000 |
| 3,3',4,4',5-PentaCB | 126 | 2000 |
| 2,2',4,4',6,6'-HexaCB | 155 | 2000 |
| 2,3,3',4,4',5-HexaCB | 156 | 2000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 2000 |
| 2,3',4,4',5,5'-HexaCB | 167 | 2000 |
| 3,3',4,4',5,5'-HexaCB | 169 | 2000 |
| 2,2',3,4',5,6,6'-HeptaCB | 188 | 2000 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 2000 |
| 2,2',3,3',5,5',6,6'-OctaCB | 202 | 2000 |
| 2,3,3',4,4',5,5',6-OctaCB | 205 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 2000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 2000 |
| DecaCB | 209 | 2000 |

CEN Method EN-1948-4 PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|---------------------------------------|
| EC-5380 | EN-1948-4 WHO PCB Calibration Series [CS1-CS6] | Set of 6 x 0.2 mL in nonane/isooctane |
| EC-5380-CS1 | EN-1948-4 WHO PCB Calibration Series [CS1] | 0.2 mL in nonane/isooctane |
| EC-5380-CS2 | EN-1948-4 WHO PCB Calibration Series [CS2] | 0.2 mL in nonane/isooctane |
| EC-5380-CS3 | EN-1948-4 WHO PCB Calibration Series [CS3] | 0.2 mL in nonane/isooctane |
| EC-5380-CS4 | EN-1948-4 WHO PCB Calibration Series [CS4] | 0.2 mL in nonane/isooctane |
| EC-5380-CS5 | EN-1948-4 WHO PCB Calibration Series [CS5] | 0.2 mL in nonane/isooctane |
| EC-5380-CS6 | EN-1948-4 WHO PCB Calibration Series [CS6] | 0.2 mL in nonane/isooctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
|--------------------------|-------|-----|-----|-----|-----|------|------|
| 3,4,4',5-TetraCB | 81 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4'-PentaCB | 105 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3',4,4',5-PentaCB | 118 | 0.6 | 6 | 60 | 300 | 1200 | 4800 |
| 2',3,4,4',5-PentaCB | 123 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 1 | 10 | 50 | 200 | 800 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 1 | 10 | 50 | 200 | 800 |

Sampling

| | | | | | | | |
|---|-----|----|----|----|----|----|----|
| 2,3,4,4'-TetraCB (¹³ C ₁₂ , 99%) | 60 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 159 | 10 | 10 | 10 | 10 | 10 | 10 |

Extraction

| | | | | | | | |
|---|-----|----|----|----|----|----|----|
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 |

Recovery

| | | | | | | | |
|---|-----|----|----|----|----|----|----|
| 2,3',4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 |

| | | | |
|-----|----------------|---------------------------------|-------|
| NEW | EC-1948-4W-KIT | EN-1948-4 WHO PCB "Starter Kit" | 1 Kit |
|-----|----------------|---------------------------------|-------|

Contains one each of the following items:

| | |
|---------|--|
| EC-5380 | EN-1948-4 WHO PCB Calibration Series [CS1-CS6] |
| EC-5370 | EN-1948-4 PCB Sampling Standard |
| EC-5372 | EN-1948-4 WHO PCB Extraction Standard |
| EC-5371 | EN-1948-4 PCB Recovery Standard |

CEN Method EN-1948-4 PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|---------------------------------------|
| EC-5385 | EN-1948-4 Marker PCB Calibration Series [CS1-CS6] | Set of 6 × 0.2 mL in nonane/isooctane |
| EC-5385-CS1 | EN-1948-4 Marker PCB Calibration Series [CS1] | 0.2 mL in nonane/isooctane |
| EC-5385-CS2 | EN-1948-4 Marker PCB Calibration Series [CS2] | 0.2 mL in nonane/isooctane |
| EC-5385-CS3 | EN-1948-4 Marker PCB Calibration Series [CS3] | 0.2 mL in nonane/isooctane |
| EC-5385-CS4 | EN-1948-4 Marker PCB Calibration Series [CS4] | 0.2 mL in nonane/isooctane |
| EC-5385-CS5 | EN-1948-4 Marker PCB Calibration Series [CS5] | 0.2 mL in nonane/isooctane |
| EC-5385-CS6 | EN-1948-4 Marker PCB Calibration Series [CS6] | 0.2 mL in nonane/isooctane |

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | | |
|--|-------|-----|-----|-----|-----|-----|------|
| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
| 2,4,4'-TriCB | 28 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| 2,2',5,5'-TetraCB | 52 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| 2,2',4,5,5'-PentaCB | 101 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.1 | 1 | 10 | 100 | 500 | 5000 |
| Sampling | | | | | | | |
| 2,3,4,4'-TetraCB ($^{13}\text{C}_{12}$, 99%) | 60 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-PentaCB ($^{13}\text{C}_{12}$, 99%) | 127 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,5,5'-HexaCB ($^{13}\text{C}_{12}$, 99%) | 159 | 10 | 10 | 10 | 10 | 10 | 10 |
| Extraction | | | | | | | |
| 2,4,4'-TriCB ($^{13}\text{C}_{12}$, 99%) | 28 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',5,5'-TetraCB ($^{13}\text{C}_{12}$, 99%) | 52 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-PentaCB ($^{13}\text{C}_{12}$, 99%) | 101 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-HexaCB ($^{13}\text{C}_{12}$, 99%) | 138 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaCB ($^{13}\text{C}_{12}$, 99%) | 153 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB ($^{13}\text{C}_{12}$, 99%) | 180 | 100 | 100 | 100 | 100 | 100 | 100 |
| Recovery | | | | | | | |
| 2,3',4',5-TetraCB ($^{13}\text{C}_{12}$, 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB ($^{13}\text{C}_{12}$, 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB ($^{13}\text{C}_{12}$, 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 |

| | | | |
|------------|----------------|------------------------------------|-------|
| NEW | EC-1948-4M-KIT | EN-1948-4 Marker PCB "Starter Kit" | 1 Kit |
|------------|----------------|------------------------------------|-------|

Contains one each of the following items:

| | |
|---------|---|
| EC-5385 | EN-1948-4 Marker PCB Calibration Series [CS1-CS6] |
| EC-5370 | EN-1948-4 PCB Sampling Standard |
| EC-5379 | EN-1948-4 Marker PCB Extraction Standard |
| EC-5371 | EN-1948-4 PCB Recovery Standard |

CEN Method EN-1948-4 PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------|--|------------------|
| EC-5370 | EN-1948-4 PCB Sampling Standard | 1.2 mL in nonane |
| EC-5370-1/10X-10 | EN-1948-4 PCB Sampling Standard (1/10 concentration) | 10 mL in nonane |

| Labeled | IUPAC | EC-5370 (ng/mL) | EC-5370-1/10X-10 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 2,3,4,4'-TetraCB (¹³ C ₁₂ , 99%) | 60 | 100 | 10 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 100 | 10 |
| 2,3,3',4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 159 | 100 | 10 |

| | | |
|------------------|--|------------------|
| EC-5372 | EN-1948-4 WHO PCB Extraction Standard | 1.2 mL in nonane |
| EC-5372-1/10X-10 | EN-1948-4 WHO PCB Extraction Standard (1/10 concentration) | 10 mL in nonane |

| Labeled | IUPAC | EC-5372 (ng/mL) | EC-5372-1/10X-10 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 100 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 100 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 100 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 100 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 100 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 100 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 100 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 100 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 100 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 100 | 10 |

| | | |
|------------------|---|----------------------|
| EC-5379 | EN-1948-4 Marker PCB Extraction Standard | 1.2 mL in nonane |
| EC-5379-5X1.2 | EN-1948-4 Marker PCB Extraction Standard | 5 × 1.2 mL in nonane |
| EC-5379-1/10X-10 | EN-1948-4 Marker PCB Extraction Standard (1/10 concentration) | 10 mL in nonane |

| Labeled | IUPAC | EC-5379 (ng/mL) | EC-5379-1/10X-10 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 1000 | 100 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 1000 | 100 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 1000 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 1000 | 100 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 1000 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | 100 |

| | | |
|---------|---------------------------------|------------------|
| EC-5371 | EN-1948-4 PCB Recovery Standard | 1.2 mL in nonane |
|---------|---------------------------------|------------------|

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 100 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 100 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 100 |

JIS PCB Methods Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|---|---------------------------------------|
| EC-5323-H | Modified JIS PCB Calibration Solutions [CS1H-CS5H] | Set of 5 × 0.2 mL in nonane/isooctane |
| EC-5323-H-E | Modified JIS PCB Extended Calibration Solutions [CS0.4H-CS6H] | Set of 7 × 0.2 mL in nonane/isooctane |
| EC-5323-CS0.4H | Modified JIS PCB Extended Calibration Solution [CS0.4H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS1H | Modified JIS PCB Extended Calibration Solution [CS1H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS2H | Modified JIS PCB Extended Calibration Solution [CS2H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS3H | Modified JIS PCB Extended Calibration Solution [CS3H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS4H | Modified JIS PCB Extended Calibration Solution [CS4H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS5H | Modified JIS PCB Extended Calibration Solution [CS5H] | 0.2 mL in nonane/isooctane |
| EC-5323-CS6H | Modified JIS PCB Extended Calibration Solution [CS6H] | 0.2 mL in nonane/isooctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS0.4H | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
|---|-------|--------|------|------|------|------|------|------|
| 3,4,4',5-TetraCB | 81 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2',3,4,4',5-PentaCB | 123 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5-PentaCB | 118 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4'-PentaCB | 105 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| Cleanup | | | | | | | | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | | |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |

JIS PCB Methods Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|---|--------------------------------------|
| EC-5360 | Modified JIS PCB Alternate A Extended Calibration Solutions [CS0.4H-CS6H] | Set of 7 x 0.2 mL in nonane/isoctane |
| EC-5360-CS0.4H | Modified JIS PCB Alternate A Extended Calibration Solution [CS0.4H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS1H | Modified JIS PCB Alternate A Extended Calibration Solution [CS1H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS2H | Modified JIS PCB Alternate A Extended Calibration Solution [CS2H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS3H | Modified JIS PCB Alternate A Extended Calibration Solution [CS3H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS4H | Modified JIS PCB Alternate A Extended Calibration Solution [CS4H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS5H | Modified JIS PCB Alternate A Extended Calibration Solution [CS5H] | 0.2 mL in nonane/isoctane |
| EC-5360-CS6H | Modified JIS PCB Alternate A Extended Calibration Solution [CS6H] | 0.2 mL in nonane/isoctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS0.4H | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
|---|-------|--------|------|------|------|------|------|------|
| 3,4,4',5-TetraCB | 81 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2',3,4,4',5-PentaCB | 123 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| Cleanup | | | | | | | | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | | |
| 2,3',4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |

JIS PCB Methods Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--------------------------------|------------------|
| EC-5324 | Modified JIS PCB Cleanup Spike | 1.2 mL in nonane |

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 50 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 50 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 50 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 50 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 50 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 50 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 50 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 100 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 50 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 50 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 50 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 50 |

| | | |
|--------------|--------------------------------|------------------|
| EC-5325 | Modified JIS PCB Syringe Spike | 1.2 mL in nonane |
| EC-5325-0.2X | Modified JIS PCB Syringe Spike | 10 mL in nonane |
| EC-5325-20X | Modified JIS PCB Syringe Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5325 (ng/mL) | EC-5325-0.2X (ng/mL) | EC-5325-20X (ng/mL) |
|---|-------|--------------------|-------------------------|------------------------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 50 | 10 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 50 | 10 | 1000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 50 | 10 | 1000 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 50 | 10 | 1000 |

| | | |
|------------------------|---------------------------------|------------------|
| EC-5326 | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |
| NEW EC-5326-20X | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5326 (ng/mL) | EC-5326-20X (ng/mL) |
|--|-------|--------------------|------------------------|
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 50 | 1000 |

JIS PCB Methods Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|---|--------------------------------------|
| EC-5418 | Modified JIS PCB Alternate B Calibration Solutions [CS1H-CS5H] | Set of 5 × 0.2 mL in nonane/isoctane |
| EC-5418-CS0.4H | Modified JIS PCB Alternate B Calibration Solution [CS0.4H] (not included with EC-5418) | 0.2 mL in nonane/isoctane |
| EC-5418-CS6H | Modified JIS PCB Alternate B Calibration Solution [CS6H] (not included with EC-5418) | 0.2 mL in nonane/isoctane |
| EC-5418-CS1H | Modified JIS PCB Alternate B Calibration Solution [CS1H] | 0.2 mL in nonane/isoctane |
| EC-5418-CS2H | Modified JIS PCB Alternate B Calibration Solution [CS2H] | 0.2 mL in nonane/isoctane |
| EC-5418-CS3H | Modified JIS PCB Alternate B Calibration Solution [CS3H] | 0.2 mL in nonane/isoctane |
| EC-5418-CS4H | Modified JIS PCB Alternate B Calibration Solution [CS4H] | 0.2 mL in nonane/isoctane |
| EC-5418-CS5H | Modified JIS PCB Alternate B Calibration Solution [CS5H] | 0.2 mL in nonane/isoctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS0.4H | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
|---|-------|--------|------|------|------|------|------|------|
| 3,4,4',5-TetraCB | 81 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2',3,4,4',5-PentaCB | 123 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| Cleanup | | | | | | | | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 8 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 8 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 8 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | | |
| 2,3',4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 4 | 10 | 10 | 10 | 10 | 10 | 10 |

JIS PCB Methods Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|------------------|
| EC-5419 | Modified JIS PCB Alternate B Cleanup Solution | 1.2 mL in nonane |

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 50 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 50 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 50 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 50 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 50 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 50 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 100 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 50 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 50 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 50 |

| | | |
|---------|--|------------------|
| EC-5420 | Modified JIS PCB Alternate B Syringe Spike | 1.2 mL in nonane |
| EC-5163 | PCB Mixture (PCB-70/111/138/170) | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5420 (ng/mL) | EC-5163 (ng/mL) |
|---|-------|--------------------|--------------------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 50 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 50 | 1000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 50 | 1000 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 50 | 1000 |

| | | |
|------------------------|---------------------------------|------------------|
| EC-5326 | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |
| NEW EC-5326-20X | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5326 (ng/mL) | EC-5326-20X (ng/mL) |
|--|-------|--------------------|------------------------|
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 50 | 1000 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | Amount |
|----------------|--|-----------------------------|
| EC-5421-H-E | DL-PCB RH12 Extended Calibration Solutions [CS0.4H-CS6H] | Set of 7 x 0.2 mL in nonane |
| EC-5421-H | DL-PCB RH12 Calibration Solutions [CS1H-CS5H] | Set of 5 x 0.2 mL in nonane |
| EC-5421-CS0.4H | DL-PCB RH12 Extended Calibration Solution [CS0.4H] | 0.2 mL in nonane |
| EC-5421-CS1H | DL-PCB RH12 Calibration Solution [CS1H] | 0.2 mL in nonane |
| EC-5421-CS2H | DL-PCB RH12 Calibration Solution [CS2H] | 0.2 mL in nonane |
| EC-5421-CS3H | DL-PCB RH12 Calibration Solution [CS3H] | 0.2 mL in nonane |
| EC-5421-CS4H | DL-PCB RH12 Calibration Solution [CS4H] | 0.2 mL in nonane |
| EC-5421-CS5H | DL-PCB RH12 Calibration Solution [CS5H] | 0.2 mL in nonane |
| EC-5421-CS6H | DL-PCB RH12 Extended Calibration Solution [CS6H] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS0.4H | CS1H | CS2H | CS3H | CS4H | CS5H | CS6H |
|---|-------|--------|------|------|------|------|------|------|
| 3,4,4',5-TetraCB | 81 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2',3,4,4',5-PentaCB | 123 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 0.25 | 1 | 5 | 20 | 100 | 500 |
| Extraction | | | | | | | | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 20 | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | | |
| 2,3',4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | | |
| 2,3,4,4'-TetraCB (¹³ C ₁₂ , 99%) | 60 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 159 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | Amount |
|------------------------|------------------------------|------------------|
| EC-5422 | DL-PCB RH12 Extraction Spike | 1.2 mL in nonane |
| NEW EC-5422-10X | DL-PCB RH12 Extraction Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5422 (ng/mL) | EC-5422-10X (ng/mL) |
|---|-------|--------------------|------------------------|
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 100 | 1000 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 100 | 1000 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 100 | 1000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 200 | 2000 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 100 | 1000 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 200 | 2000 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 100 | 1000 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 100 | 1000 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 200 | 2000 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 100 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 100 | 1000 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 100 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 100 | 1000 |

| | | |
|------------------------|---------------------------|------------------|
| EC-5423 | DL-PCB RH12 Syringe Spike | 1.2 mL in nonane |
| NEW EC-5423-10X | DL-PCB RH12 Syringe Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5423 (ng/mL) | EC-5423-10X (ng/mL) |
|---|-------|--------------------|------------------------|
| 2,2',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 100 | 1000 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 100 | 1000 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 100 | 1000 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 100 | 1000 |

| | | |
|------------------------|----------------------------|------------------|
| EC-5424 | DL-PCB RH12 Sampling Spike | 1.2 mL in nonane |
| NEW EC-5424-10X | DL-PCB RH12 Sampling Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5424 (ng/mL) | EC-5424-10X (ng/mL) |
|---|-------|--------------------|------------------------|
| 2,3,4,4'-TetraCB (¹³ C ₁₂ , 99%) | 60 | 100 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 100 | 1000 |
| 2,3,3',4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 159 | 100 | 1000 |

| | | |
|------------------------|--|------------------|
| NEW EC-5505-10X | DL-PCB RH12 Native Spike (excluding PCB-170) | 1.2 mL in nonane |
|------------------------|--|------------------|

| Labeled | IUPAC | (ng/mL) |
|--------------------------|-------|---------|
| 3,4,4',5-TetraCB | 81 | 1000 |
| 3,3',4,4'-TetraCB | 77 | 1000 |
| 2',3,4,4',5-PentaCB | 123 | 1000 |
| 2,3',4,4',5-PentaCB | 118 | 2000 |
| 2,3,4,4',5-PentaCB | 114 | 1000 |
| 2,3,3',4,4'-PentaCB | 105 | 2000 |
| 3,3',4,4',5-PentaCB | 126 | 1000 |
| 2,3',4,4',5,5'-HexaCB | 167 | 1000 |
| 2,3,3',4,4',5-HexaCB | 156 | 2000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 1000 |
| 3,3',4,4',5,5'-HexaCB | 169 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 1000 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|-----------------------------|
| EC-5396 | Co-PCB Calibration Solutions [CS1-CS6] | Set of 6 x 0.2 mL in nonane |
| EC-5396-CS1 | Co-PCB Calibration Solution [CS1] | 0.2 mL in nonane |
| EC-5396-CS2 | Co-PCB Calibration Solution [CS2] | 0.2 mL in nonane |
| EC-5396-CS3 | Co-PCB Calibration Solution [CS3] | 0.2 mL in nonane |
| EC-5396-CS4 | Co-PCB Calibration Solution [CS4] | 0.2 mL in nonane |
| EC-5396-CS5 | Co-PCB Calibration Solution [CS5] | 0.2 mL in nonane |
| EC-5396-CS6 | Co-PCB Calibration Solution [CS6] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
|---|-------|-----|-----|-----|-----|-----|-----|
| 3,4,4',5-TetraCB | 81 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 3,3',4,4'-TetraCB | 77 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 3,3',4,4',5-PentaCB | 126 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2',3,4,4',5-PentaCB | 123 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,4,4',5-PentaCB | 114 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.2 | 0.5 | 2 | 10 | 50 | 200 |
| Labeled | | | | | | | |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 | 10 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EC-5315 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solutions (low) [CS1-CS5] | Set of 5 × 0.2 mL in nonane |
| EC-5315-1 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solution [CS1] | 0.2 mL in nonane |
| EC-5315-2 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solution [CS2] | 0.2 mL in nonane |
| EC-5315-3 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solution [CS3] | 0.2 mL in nonane |
| EC-5315-4 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solution [CS4] | 0.2 mL in nonane |
| EC-5315-5 | WHO PCB and PCB-170 + 180 + Syringe PCB Calibration Solution [CS5] | 0.2 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 |
|---|-------|-----|-----|-----|-----|-----|
| 3,3',4,4'-TetraCB | 77 | 0.2 | 1 | 5 | 20 | 100 |
| 3,4,4',5-TetraCB | 81 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,4,4',5-PentaCB | 114 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 1 | 5 | 20 | 100 |
| 2',3,4,4',5-PentaCB | 123 | 0.2 | 1 | 5 | 20 | 100 |
| 3,3',4,4',5-PentaCB | 126 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.2 | 1 | 5 | 20 | 100 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.2 | 1 | 5 | 20 | 100 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.2 | 1 | 5 | 20 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.2 | 1 | 5 | 20 | 100 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.2 | 1 | 5 | 20 | 100 |
| Labeled | | | | | | |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|--------------------------------------|
| EC-5186 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Calibration Solutions [CS1-CS5] | Set of 5 × 0.2 mL in nonane/isoctane |
| EC-5186-CS1 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Solution [CS1] | 0.2 mL in nonane/isoctane |
| EC-5186-CS2 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Solution [CS2] | 0.2 mL in nonane/isoctane |
| EC-5186-CS3 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Solution [CS3] | 0.2 mL in nonane/isoctane |
| EC-5186-CS4 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Solution [CS4] | 0.2 mL in nonane/isoctane |
| EC-5186-CS5 | WHO PCB + PCB-170 + PCB-180 + Syringe PCB Solution [CS5] | 0.2 mL in nonane/isoctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 |
|---|-------|-----|-----|-----|-----|-----|
| 3,3',4,4'-TetraCB | 77 | 0.5 | 2 | 10 | 50 | 250 |
| 3,4,4',5-TetraCB | 81 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,4',5-PentaCB | 126 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.5 | 2 | 10 | 50 | 250 |
| 2',3,4,4',5-PentaCB | 123 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4',5-PentaCB | 118 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4'-PentaCB | 105 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,4,4',5-PentaCB | 114 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.5 | 2 | 10 | 50 | 250 |
| Labeled | | | | | | |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 |
| Sampling | | | | | | |
| 3,3',4,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 10 | 10 | 10 | 10 | 10 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | | Amount |
|-------------|---|-------|----------------|
| EC-4937 | WHO Coplanar and Mono-Ortho PCBs | | 3 mL in nonane |
| | Labeled | IUPAC | (ng/mL) |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 |
| | 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 |
| | 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 |
| | 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 |
| | 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 |
| | 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 |
| | 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 |
| | 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 |
| | 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 |
| | 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 |

| | | |
|---------|---|------------------|
| EC-4995 | WHO Coplanar and Mono-Ortho PCBs with PCB 170/180 | 1.2 mL in nonane |
| EC-5045 | WHO PCB + PCB-170 + PCB-180 Cleanup Standard | 1.2 mL in nonane |

| Labeled | IUPAC | EC-4995 (ng/mL) | EC-5045 (ng/mL) |
|---|-------|--------------------|--------------------|
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 | 2000 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 | 2000 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 | 2000 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 | 2000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 | 2000 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 | 2000 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 | 2000 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 | 2000 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 | 2000 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 | 2000 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 | 2000 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 1000 | 2000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | 2000 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 | 2000 |

| | | |
|-----------------|-------------------------------|------------------|
| EC-5397 | Co-PCB Syringe Spike | 1.2 mL in nonane |
| EC-5181 | PCB Syringe Spike | 1.2 mL in nonane |
| EC-5181-10X-1.2 | PCB Syringe Spike (10X stock) | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5397 (ng/mL) | EC-5181 (ng/mL) | EC-5181-10X-1.2 (ng/mL) |
|---|-------|--------------------|--------------------|----------------------------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 20 | 100 | 1000 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 20 | 100 | 1000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 20 | 100 | 1000 |
| 2,2',3,3',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 20 | — | — |

| | | |
|-----------------|---------------------------------|------------------|
| EC-5326 | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |
| NEW EC-5326-20X | Modified JIS PCB Sampling Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5326 (ng/mL) | EC-5326-20X (ng/mL) |
|--|-------|--------------------|------------------------|
| 3,3',4,5,5'-TetraCB (¹³ C ₁₂ , 99%) | 79 | 50 | 1000 |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | IUPAC | (ng/mL) | Amount |
|-------------|---|-------|---------|----------------|
| EC-4070 | Coplanar PCB Mixture | | | 3 mL in nonane |
| | Labeled | | | |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 5000 | |
| | 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 5000 | |
| | 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 5000 | |
| EC-4187 | Coplanar PCB Mixture | | | 3 mL in nonane |
| | Labeled | | | |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 | |
| | 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 | |
| | 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 | |
| | 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 | |
| EC-4188 | Mono-Ortho PCB Mixture – *High Purity | | | 3 mL in nonane |
| | Labeled | | | |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 | |
| | 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 | |
| | 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 | |
| | 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 | |
| | 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 | |
| | 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 | |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 | |
| | 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 | |
| EC-4938 | PCB Mixture-A | | | 3 mL in nonane |
| | Labeled | | | |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 | |
| | 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 | |
| | 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 | |
| | 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 | |
| | 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 | |
| | 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | |

WHO "Dioxin-Like" PCB Mixtures

| Catalog No. | Compound | | Amount |
|--------------------------|--|--------------------|------------------------|
| EC-4935 | WHO Coplanar and Mono-Ortho PCBs | | 1.2 mL in isoctane |
| EC-4935-A | WHO Coplanar and Mono-Ortho PCBs | | 3 mL in isoctane |
| EC-4935-B | WHO Coplanar and Mono-Ortho PCBs | | 1.2 mL in isoctane |
| NEW EC-5559 | DL-PCB Native Plus Mixture | | 1.2 mL in isoctane |
| | | | |
| Unlabeled | IUPAC | EC-4935 (ng/mL) | EC-4935-A (ng/mL) |
| 3,3',4,4'-TetraCB | 77 | 2000 | 1000 |
| 3,4,4',5-TetraCB | 81 | 2000 | 1000 |
| 2,3,3',4,4'-PentaCB | 105 | 2000 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 2000 | 1000 |
| 2,3',4,4',5-PentaCB | 118 | 2000 | 1000 |
| 2',3,4,4',5-PentaCB | 123 | 2000 | 1000 |
| 3,3',4,4',5-PentaCB | 126 | 2000 | 1000 |
| 2,3,3',4,4',5-HexaCB | 156 | 2000 | 1000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 2000 | 1000 |
| 2,3',4,4',5,5'-HexaCB | 167 | 2000 | 1000 |
| 3,3',4,4',5,5'-HexaCB | 169 | 2000 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 2000 | 1000 |
| | | | |
| EC-5000 | WHO Coplanar and Mono-Ortho PCBs and 170/180 | | 1.2 mL in isoctane |
| | | | |
| Unlabeled | IUPAC | (ng/mL) | |
| 3,3',4,4'-TetraCB | 77 | 2000 | |
| 3,4,4',5-TetraCB | 81 | 2000 | |
| 2,3,3',4,4'-PentaCB | 105 | 2000 | |
| 2,3,4,4',5-PentaCB | 114 | 2000 | |
| 2,3',4,4',5-PentaCB | 118 | 2000 | |
| 2',3,4,4',5-PentaCB | 123 | 2000 | |
| 3,3',4,4',5-PentaCB | 126 | 2000 | |
| 2,3,3',4,4',5-HexaCB | 156 | 2000 | |
| 2,3,3',4,4',5'-HexaCB | 157 | 2000 | |
| 2,3',4,4',5,5'-HexaCB | 167 | 2000 | |
| 3,3',4,4',5,5'-HexaCB | 169 | 2000 | |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 2000 | |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 2000 | |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 2000 | |
| | | | |
| EC-4986 | Non-Ortho Native PCB Solution | | 1.2 mL in isoctane |
| | | | |
| Unlabeled | IUPAC | (ng/mL) | |
| 3,3',4,4'-TetraCB | 77 | 10,000 | |
| 3,4,4',5-TetraCB | 81 | 10,000 | |
| 3,3',4,4',5-PentaCB | 126 | 10,000 | |
| 3,3',4,4',5,5'-HexaCB | 169 | 10,000 | |
| | | | |
| EC-4987 | Mono-Ortho Native PCB Solution | | 1.2 mL in isoctane |
| EC-4987/100 | Mono-Ortho Native PCB Solution | | 100 µL in isoctane |
| | | | |
| Unlabeled | IUPAC | EC-4987 (ng/mL) | EC-4987/100 (ng/mL) |
| 2,3,3',4,4'-PentaCB | 105 | 10,000 | 100 |
| 2,3,4,4',5-PentaCB | 114 | 10,000 | 100 |
| 2,3',4,4',5-PentaCB | 118 | 10,000 | 100 |
| 2',3,4,4',5-PentaCB | 123 | 10,000 | 100 |
| 2,3,3',4,4',5-HexaCB | 156 | 10,000 | 100 |
| 2,3,3',4,4',5'-HexaCB | 157 | 10,000 | 100 |
| 2,3',4,4',5,5'-HexaCB | 167 | 10,000 | 100 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 10,000 | 100 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|--------------------------------------|
| EC-5414 | Mono-Deca Plus Predominant PCB Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane/isoctane |
| EC-5414-CS1 | Mono-Deca Plus Predominant PCB Calibration Solution [CS1] | 0.2 mL in nonane/isoctane |
| EC-5414-CS2 | Mono-Deca Plus Predominant PCB Calibration Solution [CS2] | 0.2 mL in nonane/isoctane |
| EC-5414-CS3 | Mono-Deca Plus Predominant PCB Calibration Solution [CS3] | 0.2 mL in nonane/isoctane |
| EC-5414-CS4 | Mono-Deca Plus Predominant PCB Calibration Solution [CS4] | 0.2 mL in nonane/isoctane |
| EC-5414-CS5 | Mono-Deca Plus Predominant PCB Calibration Solution [CS5] | 0.2 mL in nonane/isoctane |

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | |
|---|-------|-----|-----|-----|-----|------|
| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 |
| 4-MonoCB | 3 | 4 | 20 | 100 | 500 | 2000 |
| 2,4'-DiCB | 8 | 4 | 20 | 100 | 500 | 2000 |
| 2,4,4'-TriCB | 28 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',5,5'-TetraCB | 52 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',4,5,5'-PentaCB | 101 | 2 | 10 | 50 | 250 | 1000 |
| 2,3',4,4',5-PentaCB | 118 | 2 | 10 | 50 | 250 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 2 | 10 | 50 | 250 | 1000 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 4 | 20 | 100 | 500 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 4 | 20 | 100 | 500 | 2000 |
| DecaCB | 209 | 4 | 20 | 100 | 500 | 2000 |
| Labeled | | | | | | |
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 100 | 100 | 100 | 100 | 100 |
| 2,4'-DiCB (¹³ C ₁₂ , 99%) | 8 | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 | 100 | 100 | 100 | 100 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 100 | 100 | 100 | 100 | 100 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 100 | 100 | 100 | 100 | 100 |
| Syringe | | | | | | |
| 2,4',6-TriCB (¹³ C ₁₂ , 99%) | 32 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 100 | 100 | 100 | 100 | 100 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|----------------------|--|------------------|
| EC-5411 | Predominant Mono-Deca PCB Spike | 1.2 mL in nonane |
| NEW EC-5411-A | Modified Predominant Mono-Deca PCB Spike | 1.2 mL in nonane |

| Labeled | IUPAC | EC-5411 (ng/mL) | EC-5411-A (ng/mL) |
|---|-------|--------------------|----------------------|
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 2000 | 2000 |
| 2,4'-DiCB (¹³ C ₁₂ , 99%) | 8 | 2000 | 2000 |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 2000 | 1000 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 2000 | 1000 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 2000 | 1000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 2000 | 1000 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 2000 | 1000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 2000 | 1000 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 2000 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 2000 | 1000 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 2000 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 2000 | 2000 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 2000 | 2000 |

| | | |
|---------|--|------------------|
| EC-5415 | Mono-Deca Plus Predominant PCB Syringe Spike | 1.2 mL in nonane |
|---------|--|------------------|

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,4',6-TriCB (¹³ C ₁₂ , 99%) | 32 | 2000 |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 2000 |
| 3,3',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 127 | 2000 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 2000 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 2000 |

| | | |
|---------|--|--------------------|
| EC-5412 | Native Mono-Deca Plus Predominant Spike PCBs | 1.2 mL in isoctane |
|---------|--|--------------------|

| Unlabeled | IUPAC | (ng/mL) |
|------------------------------|-------|---------|
| 4-MonoCB | 3 | 2000 |
| 2,4'-DiCB | 8 | 2000 |
| 2,4,4'-TriCB | 28 | 2000 |
| 2,2',5,5'-TetraCB | 52 | 2000 |
| 2,2',4,5,5'-PentaCB | 101 | 2000 |
| 2,3',4,4',5-PentaCB | 118 | 2000 |
| 2,3,4,4',5-PentaCB | 114 | 2000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 2000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 2000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 2000 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 2000 |
| DecaCB | 209 | 2000 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|---------------------------|--|---------------------------------------|
| <i>NEW</i> EC-5448-HR | Rapid PCB Screening Calibration Solutions [CS0.02, 0.05, 0.1, 0.5] | Set of 4 × 0.2 mL in nonane/isooctane |
| <i>NEW</i> EC-5448-CS0.02 | HRMS Rapid PCB Screening Calibration Solution [CS0.02] | 0.2 mL in nonane/isooctane |
| <i>NEW</i> EC-5448-CS0.05 | HRMS Rapid PCB Screening Calibration Solution [CS0.05] | 0.2 mL in nonane/isooctane |
| <i>NEW</i> EC-5448-CS0.1 | HRMS Rapid PCB Screening Calibration Solution [CS0.1] | 0.2 mL in nonane/isooctane |
| <i>NEW</i> EC-5448-CS0.5 | HRMS Rapid PCB Screening Calibration Solution [CS0.5] | 0.2 mL in nonane/isooctane |
| EC-5448 | Rapid PCB Screening Calibration Solutions [CS1-CS4] | Set of 4 × 0.2 mL in nonane/isooctane |
| EC-5448-CS1 | Rapid PCB Screening Calibration Solution [CS1] | 0.2 mL in nonane/isooctane |
| EC-5448-CS2 | Rapid PCB Screening Calibration Solution [CS2] | 0.2 mL in nonane/isooctane |
| EC-5448-CS3 | Rapid PCB Screening Calibration Solution [CS3] | 0.2 mL in nonane/isooctane |
| EC-5448-CS4 | Rapid PCB Screening Calibration Solution [CS4] | 0.2 mL in nonane/isooctane |

| All concentrations are in ng/mL (ppb) | | | | | | | | | |
|---|-------|--------|--------|-------|-------|-----|-----|-----|-----|
| Unlabeled | IUPAC | CS0.02 | CS0.05 | CS0.1 | CS0.5 | CS1 | CS2 | CS3 | CS4 |
| 2,4,4'-TriCB | 28 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',5-TriCB | 18 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',3,5'-TetraCB | 44 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,3',4',5-TetraCB | 70 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',5,5'-TetraCB | 52 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',4,5,5'-PentaCB | 101 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,3,3',4',6-PentaCB | 110 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,3',4,4',5-PentaCB | 118 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',3,4',5',6-HexaCB | 149 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',3,4,4',5'-HexaCB | 138 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',4,4',5,5'-HexaCB | 153 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| 2,2',3,4',5,5'-HeptaCB | 187 | 0.02 | 0.05 | 0.1 | 0.5 | 1 | 5 | 10 | 50 |
| Extraction | | | | | | | | | |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| Syringe | | | | | | | | | |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 141 | 10 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|------------------|---|----------------------|
| EC-5379 | EN-1948-4 Marker PCB Extraction Standard | 1.2 mL in nonane |
| EC-5379-5X1.2 | EN-1948-4 Marker PCB Extraction Standard | 5 × 1.2 mL in nonane |
| EC-5379-1/10X-10 | EN-1948-4 Marker PCB Extraction Standard (1/10 concentration) | 10 mL in nonane |

| Labeled | IUPAC | EC-5379 (ng/mL) | EC-5379-1/10X-10 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 1000 | 100 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 1000 | 100 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 1000 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 1000 | 100 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 1000 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | 100 |

| | | |
|---------|-----------------------------------|------------------|
| EC-5450 | Rapid PCB Screening Syringe Spike | 1.2 mL in nonane |
|---------|-----------------------------------|------------------|

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 2000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 2000 |
| 2,2',3,4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 141 | 2000 |

| | | |
|---------|---|--------------------|
| EC-5453 | Rapid PCB Screening Native PAR Solution | 0.5 mL in isoctane |
|---------|---|--------------------|

| Unlabeled | IUPAC | (ng/mL) |
|--------------------------|-------|---------|
| 2,4,4'-TriCB | 28 | 1000 |
| 2,2',5-TriCB | 18 | 1000 |
| 2,2',3,5'-TetraCB | 44 | 1000 |
| 2,3',4',5-TetraCB | 70 | 1000 |
| 2,2',5,5'-TetraCB | 52 | 1000 |
| 2,2',4,5,5'-PentaCB | 101 | 1000 |
| 2,3,3',4',6-PentaCB | 110 | 1000 |
| 2,3',4,4',5-PentaCB | 118 | 1000 |
| 2,2',3,4',5',6-HexaCB | 149 | 1000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 1000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 1000 |
| 2,2',3,4',5,5',6-HeptaCB | 187 | 1000 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|-----------------------------|-------------------------------------|------------------|
| <i>NEW</i> EC-5375 | Marker-7 PCB Mixture (with PCB-118) | 1.2 mL in nonane |
| <i>NEW</i> EC-5375-1/10X-10 | Marker-7 PCB Mixture (with PCB-118) | 10 mL in nonane |

| Labeled | IUPAC | EC-5375 (ng/mL) | EC-5375-1/10X-10 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 1000 | 100 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 1000 | 100 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 1000 | 100 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 | 100 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 1000 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 1000 | 100 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 | 100 |

| | | |
|---------|-------------|----------------|
| EC-4058 | PCB Mixture | 3 mL in nonane |
|---------|-------------|----------------|

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 5000 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 5000 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 5000 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 5000 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 5000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 5000 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 5000 |

| | | |
|-----------|-----------------------|----------------|
| EC-4189-A | Mono-Deca PCB Mixture | 3 mL in nonane |
|-----------|-----------------------|----------------|

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 1000 |
| 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 1000 |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 1000 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 1000 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 1000 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 1000 |
| 2,2',3,3',4,5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 208 | 1000 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 1000 |

"Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|--------------------|-------------------------------------|--------------------|
| NEW EC-5495 | Marker-7 PCB Mixture (with PCB-118) | 1.2 mL in isoctane |
| | Unlabeled | IUPAC (ng/mL) |
| | 2,4,4'-TriCB | 28 1000 |
| | 2,2',5,5'-TetraCB | 52 1000 |
| | 2,2',4,5,5'-PentaCB | 101 1000 |
| | 2,2',4,4',5-PentaCB | 118 1000 |
| | 2,2',4,4',5,5'-HexaCB | 153 1000 |
| | 2,2',3,4,4',5'-HexaCB | 138 1000 |
| | 2,2',3,4,4',5,5'-HeptaCB | 180 1000 |
| EC-5179 | PCB Mixture | 3 mL in isoctane |
| | Unlabeled | IUPAC (ng/mL) |
| | 2,4,4'-TriCB | 28 5000 |
| | 2,2',5,5'-TetraCB | 52 5000 |
| | 2,2',4,5,5'-PentaCB | 101 5000 |
| | 2,2',3,4,4',5'-HexaCB | 138 5000 |
| | 2,2',4,4',5,5'-HexaCB | 153 5000 |
| | 2,2',3,4,4',5,5'-HeptaCB | 180 5000 |
| | DecaCB | 209 5000 |
| NEW EC-5502 | UNEP PCB Working Solution 1 | 3 mL in isoctane |
| | Unlabeled | IUPAC (ng/mL) |
| | 2,4,4'-TriCB | 28 2.0 |
| | 2,2',5,5'-TetraCB | 52 2.5 |
| | 2,2',4,5,5'-PentaCB | 101 4.0 |
| | 2,2',3,4,4',5'-HexaCB | 138 4.0 |
| | 2,2',4,4',5,5'-HexaCB | 153 5.0 |
| | 2,2',3,4,4',5,5'-HeptaCB | 180 6.0 |

See also **EN-1948-4 PCB Standard Mixtures**, pp. 86-88.

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|------------------------|---|-------------------------------|
| <i>NEW</i> EC-5518 | Comprehensive PCB Calibration Solutions [CS1-CS5] | 5 × 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS1 | Comprehensive PCB Calibration Solution [CS1] | 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS2 | Comprehensive PCB Calibration Solution [CS2] | 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS3 | Comprehensive PCB Calibration Solution [CS3] | 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS4 | Comprehensive PCB Calibration Solution [CS4] | 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS5 | Comprehensive PCB Calibration Solution [CS5] | 0.2 mL in nonane/isoctane |
| <i>NEW</i> EC-5518-CS6 | Comprehensive PCB Calibration Solution [CS6] | 0.2 mL in nonane/isoctane |

All concentrations are in ng/mL (ppb)

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
|-----------------------|-------|-----|-----|-----|-----|-----|-----|
| 2-MonoCB | 1 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 4-MonoCB | 3 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 2,2'-DiCB | 4 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 2,4'-DiCB | 8 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 2,5-DiCB | 9 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 2,6-DiCB | 10 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 3,3'-DiCB | 11 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 3,4-DiCB | 12 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 4,4'-DiCB | 15 | 0.2 | 1 | 4 | 20 | 100 | 500 |
| 2,2',5-TriCB | 18 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',6-TriCB | 19 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,4,4'-TriCB | 28 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,4',5-TriCB | 31 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2',3,4-TriCB | 33 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4-TriCB | 35 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,4,4'-TriCB | 37 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,4,5-TriCB | 38 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,5'-TetraCB | 44 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,5'-TetraCB | 49 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',5,5'-TetraCB | 52 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',6,6'-TetraCB | 54 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',5-TetraCB | 57 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4'-TetraCB | 66 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4',5-TetraCB | 70 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,4,4',5-TetraCB | 74 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,4'-TetraCB | 77 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,5-TetraCB | 78 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,5'-TetraCB | 79 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,4,4',5-TetraCB | 81 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4,5'-PentaCB | 87 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,5',6-PentaCB | 95 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,4',5-PentaCB | 99 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,5,5'-PentaCB | 101 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,6,6'-PentaCB | 104 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4'-PentaCB | 105 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4',6-PentaCB | 110 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',5,5'-PentaCB | 111 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,4,4',5-PentaCB | 114 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4',5-PentaCB | 118 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4',5'-PentaCB | 123 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,4',5-PentaCB | 126 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4,4',5'-HexaCB | 138 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4',5',6-HexaCB | 149 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,4',5,5'-HexaCB | 153 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',4,4',6,6'-HexaCB | 155 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |

(continued on next page)

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

(continued from previous page)

| | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
|---|-------|-----|-----|-----|-----|-----|-----|
| Unlabeled | | | | | | | |
| 2,3,3',4',5,5'-HexaCB | 162 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 3,3',4,4',5,5'-HexaCB | 169 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,5,6'-HeptaCB | 174 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',5,5',6-HeptaCB | 178 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4',5,5',6-HeptaCB | 187 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4',5,6,6'-HeptaCB | 188 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,4',5,6-OctaCB | 195 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,5,6,6'-OctaCB | 200 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',5,5',6,6'-OctaCB | 202 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,4,4',5,5',6-OctaCB | 203 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,3,3',4,4',5,5',6-OctaCB | 205 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| DecaCB | 209 | 0.1 | 0.5 | 2 | 10 | 50 | 250 |
| Toxics/LOC/Window Defining | | | | | | | |
| 2-MonoCB (¹³ C ₁₂ , 99%) | 1 | 10 | 10 | 10 | 10 | 10 | 10 |
| 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2'-DiCB (¹³ C ₁₂ , 99%) | 4 | 10 | 10 | 10 | 10 | 10 | 10 |
| 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',6-TriCB (¹³ C ₁₂ , 99%) | 19 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,4,4'-TriCB (¹³ C ₁₂ , 99%) | 37 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',6,6'-TetraCB (¹³ C ₁₂ , 99%) | 54 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',4,6,6'-PentaCB (¹³ C ₁₂ , 99%) | 104 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5'-PentaCB (¹³ C ₁₂ , 99%) | 123 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',4,4',6,6'-HexaCB (¹³ C ₁₂ , 99%) | 155 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 10 | 10 | 10 | 10 | 10 | 10 |
| 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4',5,6,6'-HeptaCB (¹³ C ₁₂ , 99%) | 188 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6,6'-OctaCB (¹³ C ₁₂ , 99%) | 202 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',4,4',5,5',6-OctaCB (¹³ C ₁₂ , 99%) | 205 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 10 | 10 | 10 | 10 | 10 | 10 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 10 | 10 | 10 | 10 | 10 | 10 |
| Cleanup | | | | | | | |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10 | 10 | 10 | 10 | 10 | 10 |
| Injection Internal | | | | | | | |
| 2,5-DiCB (¹³ C ₁₂ , 99%) | 9 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 10 | 10 | 10 | 10 | 10 | 10 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 10 | 10 | 10 | 10 | 10 | 10 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | IUPAC | Amount (ng/mL) |
|-------------|---|-------|-------------------|
| EC-4977 | Method 1668A/B/C Labeled Toxics/LOC/Window Defining Solution | | 1.2 mL in nonane |
| EC-4977-5 | Method 1668A/B/C Labeled Toxics/LOC/Window Defining Solution | | 5 mL in nonane |
| | Labeled | | |
| | 2-MonoCB (¹³ C ₁₂ , 99%) | 1 | 1000 |
| | 4-MonoCB (¹³ C ₁₂ , 99%) | 3 | 1000 |
| | 2,2'-DiCB (¹³ C ₁₂ , 99%) | 4 | 1000 |
| | 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 1000 |
| | 2,2',6-TriCB (¹³ C ₁₂ , 99%) | 19 | 1000 |
| | 3,4,4'-TriCB (¹³ C ₁₂ , 99%) | 37 | 1000 |
| | 2,2',6,6'-TetraCB (¹³ C ₁₂ , 99%) | 54 | 1000 |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 1000 |
| | 3,4,4',5-TetraCB (¹³ C ₁₂ , 99%) | 81 | 1000 |
| | 2,2',4,6,6'-PentaCB (¹³ C ₁₂ , 99%) | 104 | 1000 |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 1000 |
| | 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 1000 |
| | 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 1000 |
| | 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 1000 |
| | 3,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 126 | 1000 |
| | 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 155 | 1000 |
| | 2,2',4,4',6,6'-HexaCB (¹³ C ₁₂ , 99%) | 156 | 1000 |
| | 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 1000 |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 1000 |
| | 3,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 169 | 1000 |
| | 2,2',3,4',5,6,6'-HeptaCB (¹³ C ₁₂ , 99%) | 188 | 1000 |
| | 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 1000 |
| | 2,2',3,3',5,5',6,6'-OctaCB (¹³ C ₁₂ , 99%) | 202 | 1000 |
| | 2,3,3',4,4',5,5',6-OctaCB (¹³ C ₁₂ , 99%) | 205 | 1000 |
| | 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 1000 |
| | 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 1000 |
| | DecaCB (¹³ C ₁₂ , 99%) | 209 | 1000 |

| | | |
|---------|---|------------------|
| EC-4978 | Method 1668A/B/C Labeled Cleanup Standard Solution | 1.2 mL in nonane |
| | Labeled | |
| | 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 |
| | 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 |
| | 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 |
| | | 1000 |
| | | 1000 |
| | | 1000 |

| | | |
|---------|---|------------------|
| EC-4979 | Method 1668A/B/C Labeled Injection Internal Standard Solution | 1.2 mL in nonane |
| | Labeled | |
| | 2,5-DiCB (¹³ C ₁₂ , 99%) | 9 |
| | 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 |
| | 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 |
| | 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 |
| | 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 |
| | | 5000 |
| | | 5000 |
| | | 5000 |
| | | 5000 |
| | | 5000 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | | Amount |
|------------------------------|----------------------------------|-------|--------------------|
| EC-5433 | Comprehensive Native PCB Mixture | | 1.2 mL in isoctane |
| | Unlabeled | IUPAC | (ng/mL) |
| 2-MonoCB | | 1 | 2000 |
| 4-MonoCB | | 3 | 2000 |
| 2,2'-DiCB | | 4 | 2000 |
| 2,4'-DiCB | | 8 | 2000 |
| 2,5-DiCB | | 9 | 2000 |
| 2,6-DiCB | | 10 | 2000 |
| 3,3'-DiCB | | 11 | 2000 |
| 3,4-DiCB | | 12 | 2000 |
| 4,4'-DiCB | | 15 | 2000 |
| 2,2',5-TriCB | | 18 | 1000 |
| 2,2',6-TriCB | | 19 | 1000 |
| 2,4,4'-TriCB | | 28 | 1000 |
| 2,4',5-TriCB | | 31 | 1000 |
| 2',3,4-TriCB | | 33 | 1000 |
| 3,3',4-TriCB | | 35 | 1000 |
| 3,4,4'-TriCB | | 37 | 1000 |
| 3,4,5-TriCB | | 38 | 1000 |
| 2,2',3,5'-TetraCB | | 44 | 1000 |
| 2,2',4,5'-TetraCB | | 49 | 1000 |
| 2,2',5,5'-TetraCB | | 52 | 1000 |
| 2,2',6,6'-TetraCB | | 54 | 1000 |
| 2,3,3',5-TetraCB | | 57 | 1000 |
| 2,2',4,4'-TetraCB | | 66 | 1000 |
| 2,3',4',5-TetraCB | | 70 | 1000 |
| 2,4,4',5-TetraCB | | 74 | 1000 |
| 3,3',4,4'-TetraCB | | 77 | 1000 |
| 3,3',4,5-TetraCB | | 78 | 1000 |
| 3,3',4,5'-TetraCB | | 79 | 1000 |
| 3,4,4',5-TetraCB | | 81 | 1000 |
| 2,2',3,4,5'-PentaCB | | 87 | 1000 |
| 2,2',3,5,6-PentaCB | | 95 | 1000 |
| 2,2',4,4',5-PentaCB | | 99 | 1000 |
| 2,2',4,5,5'-PentaCB | | 101 | 1000 |
| | Unlabeled | IUPAC | (ng/mL) |
| 2,2',4,6,6'-PentaCB | | 104 | 1000 |
| 2,3,3',4,4'-PentaCB | | 105 | 1000 |
| 2,3,3',4',6-PentaCB | | 110 | 1000 |
| 2,3,3',5,5'-PentaCB | | 111 | 1000 |
| 2,3,4,4',5-PentaCB | | 114 | 1000 |
| 2,3',4,4',5-PentaCB | | 118 | 1000 |
| 2',3,4,4',5-PentaCB | | 123 | 1000 |
| 3,3',4,4',5-PentaCB | | 126 | 1000 |
| 2,2',3,4,4',5-HexaCB | | 138 | 1000 |
| 2,2',3,4',5',6-HexaCB | | 149 | 1000 |
| 2,2',4,4',5,5'-HexaCB | | 153 | 1000 |
| 2,2',4,4',6,6'-HexaCB | | 155 | 1000 |
| 2,3,3',4,4',5-HexaCB | | 156 | 1000 |
| 2,3,3',4,4',5'-HexaCB | | 157 | 1000 |
| 2,3,3',4',5,5'-HexaCB | | 162 | 1000 |
| 2,3',4,4',5,5'-HexaCB | | 167 | 1000 |
| 3,3',4,4',5,5'-HexaCB | | 169 | 1000 |
| 2,2',3,3',4,4',5-HeptaCB | | 170 | 1000 |
| 2,2',3,3',4,5,6'-HeptaCB | | 174 | 1000 |
| 2,2',3,3',5,5',6-HeptaCB | | 178 | 1000 |
| 2,2',3,4,4',5,5'-HeptaCB | | 180 | 1000 |
| 2,2',3,4',5,5',6-HeptaCB | | 187 | 1000 |
| 2,2',3,4',5,6,6'-HeptaCB | | 188 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB | | 189 | 1000 |
| 2,2',3,3',4,4',5,5'-OctaCB | | 194 | 1000 |
| 2,2',3,3',4,4',5,6-OctaCB | | 195 | 1000 |
| 2,2',3,3',4,5,6,6'-OctaCB | | 200 | 1000 |
| 2,2',3,3',5,5',6,6'-OctaCB | | 202 | 1000 |
| 2,2',3,4,4',5,5',6-OctaCB | | 203 | 1000 |
| 2,3,3',4,4',5,5',6-OctaCB | | 205 | 1000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | | 206 | 1000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | | 208 | 1000 |
| DecaCB | | 209 | 1000 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount | | | | | | | | | |
|--|---|------------------------------|------|-----|------|-----|------|-----|-----|------|------|
| EC-5366 | CDC PCB Calibration Solutions [CS1-CS10] (NOTE: Individual calibration standards are available upon request) | Set of 10 × 0.5 mL in nonane | | | | | | | | | |
| All concentrations are in ng/mL (ppb) | | | | | | | | | | | |
| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 | CS10 |
| 2,2',5-TriCB | 18 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,4,4'-TriCB | 28 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,5'-TetraCB | 44 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',4,5'-TetraCB | 49 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',5,5'-TetraCB | 52 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3',4,4'-TetraCB | 66 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,4,4',5-TetraCB | 74 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4,5'-PentaCB | 87 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',4,4',5-PentaCB | 99 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',4,5,5'-PentaCB | 101 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3,3',4',6-PentaCB | 110 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3,4,4',5-PentaCB | 114 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | 3000 | 7500 |
| 2',3,4,4',5-PentaCB | 123 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4,4'-HexaCB | 128 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4,4',5-HexaCB | 138 | 0.1 | 0.25 | 0.5 | 1.25 | 5 | 37.5 | 50 | 250 | 1500 | 3750 |
| 2,2',3,4',5,5'-HexaCB | 146 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4',5',6-HexaCB | 149 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,5,5',6-HexaCB | 151 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',4,4',5,5'-HexaCB | 153 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | 3000 | 7500 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3,3',4,4',5'-HexaCB | 157 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,3,3',4,4',6-HexaCB | 158 | 0.1 | 0.25 | 0.5 | 1.25 | 5 | 37.5 | 50 | 250 | 1500 | 3750 |
| 2,3',4,4',5,5'-HexaCB | 167 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | 3000 | 7500 |
| 2,2',3,3',4,5,5'-HeptaCB | 172 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4',5,6-HeptaCB | 177 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',5,5',6-HeptaCB | 178 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | 3000 | 7500 |
| 2,2',3,4,4',5',6-HeptaCB | 183 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4',5,5',6-HeptaCB | 187 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | 3000 | 7500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4,4',5,6-OctaCB | 195 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,3',4,4',5',6-OctaCB | 196 | 0.1 | 0.25 | 0.5 | 1.25 | 5 | 37.5 | 50 | 250 | | |
| 2,2',3,3',4,5,5',6'-OctaCB | 201 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| 2,2',3,4,4',5,5',6-OctaCB | 203 | 0.1 | 0.25 | 0.5 | 1.25 | 5 | 37.5 | 50 | 250 | | |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| DecaCB | 209 | 0.2 | 0.5 | 1 | 2.5 | 10 | 75 | 100 | 500 | | |
| Labeled | | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 | CS10 |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |

(continued on next page)

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

(continued from previous page)

| Labeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 | CS10 |
|---|-------|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Recovery | | | | | | | | | | | |
| 1,2,3,4-TCDD (¹³ C ₆ , 99%) | | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |

| | | |
|-------------------------|--------------------------|-----------------------|
| EC-5367 | CDC PCB Spiking Standard | 10 mL in methanol |
| NEW EC-5367-5X10 | CDC PCB Spiking Standard | 5 × 10 mL in methanol |

| Labeled | IUPAC | (ng/mL) |
|---|-------|---------|
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 7.5 |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 7.5 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 7.5 |
| 2',3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 123 | 7.5 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 7.5 |
| 2,3,4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 114 | 7.5 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 7.5 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 7.5 |
| 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 7.5 |
| 2,2',3,4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 138 | 7.5 |
| 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 128 | 7.5 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 7.5 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 7.5 |
| 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 7.5 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 7.5 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 7.5 |
| 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 7.5 |
| 2,2',3,3',4,4',5,5'-OctaCB (¹³ C ₁₂ , 99%) | 194 | 7.5 |
| 2,2',3,3',4,4',5,5',6-NonaCB (¹³ C ₁₂ , 99%) | 206 | 7.5 |
| DecaCB (¹³ C ₁₂ , 99%) | 209 | 7.5 |

| | | |
|-----------------------------|---|--|
| ES-5321 | Multi-Analyte Recovery Spiking Standard | 10 mL in 88% hexane/ 2% dodecane/10% nonane |
| NEW ES-5321-200X-1.2 | Multi-Analyte Recovery Spiking Standard | 1.2 mL in nonane |

| Labeled | IUPAC | ES-5321 (ng/mL) | ES-5321-200X-1.2 (ng/mL) |
|---|-------|--------------------|-----------------------------|
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 2.5 | 500 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 10.0 | 2000 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 7.5 | 1500 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 | 7.5 | 1500 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount |
|------------------------------|--|------------------|
| EC-5085 | Toxic and Predominant PCB PAR Solution (does not include PCB-114 or PCB-123) | 1.2 mL in nonane |
| Unlabeled | | |
| 2,2',5-TriCB | IUPAC | (ng/mL) |
| 2,2',5-TriCB | 18 | 250 |
| 2,4,4'-TriCB | 28 | 250 |
| 2,2',3,5'-TetraCB | 44 | 250 |
| 2,2',4,5'-TetraCB | 49 | 250 |
| 2,2',5,5'-TetraCB | 52 | 250 |
| 2,3',4,4'-TetraCB | 66 | 250 |
| 2,4,4',5-TetraCB | 74 | 250 |
| 2,2',3,4,5'-PentaCB | 87 | 250 |
| 2,2',4,4',5-PentaCB | 99 | 250 |
| 2,2',4,5,5'-PentaCB | 101 | 250 |
| 2,3,3',4,4'-PentaCB | 105 | 250 |
| 2,3,3',4',6-PentaCB | 110 | 250 |
| 2,3',4,4',5-PentaCB | 118 | 250 |
| 2,2',3,3',4,4'-HexaCB | 128 | 250 |
| 2,2',3,4,4',5'-HexaCB | 138 | 250 |
| 2,2',3,4',5,5'-HexaCB | 146 | 250 |
| 2,2',3,4',5',6-HexaCB | 149 | 250 |
| 2,2',3,5,5',6-HexaCB | 151 | 250 |
| 2,2',4,4',5,5'-HexaCB | 153 | 250 |
| 2,3,3',4,4',5-HexaCB | 156 | 250 |
| 2,3,3',4,4',5'-HexaCB | 157 | 250 |
| 2,3,3',4,4',6-HexaCB | 158 | 250 |
| 2,3',4,4',5,5'-HexaCB | 167 | 250 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 250 |
| 2,2',3,3',4,5,5'-HeptaCB | 172 | 250 |
| 2,2',3,3',4',5,6-HeptaCB | 177 | 250 |
| 2,2',3,3',5,5',6-HeptaCB | 178 | 250 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 250 |
| 2,2',3,4,4',5',6-HeptaCB | 183 | 250 |
| 2,2',3,4',5,5',6-HeptaCB | 187 | 250 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 250 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 250 |
| 2,2',3,3',4,4',5,6-OctaCB | 195 | 250 |
| 2,2',3,3',4,4',5',6-OctaCB | 196 | 250 |
| 2,2',3,3',4,5,5',6-OctaCB | 201 | 250 |
| 2,2',3,4,4',5,5',6-OctaCB | 203 | 250 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 250 |
| DecaCB | | 209 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | Amount | | | | | |
|---|----------|--------------------------------|-----|-----|-----|-----|------|
| NEW | EC-5531 | 6 × 0.2 mL in nonane/isooctane | | | | | |
| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | | |
| Unlabeled | IUPAC | CS0.2 | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,2',5,5'-TetraCB | 52 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,3',4,4'-TetraCB | 66 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,4,4',5-TetraCB | 74 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,4,5'-PentaCB | 87 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',4,4',5-PentaCB | 99 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',4,5,5'-PentaCB | 101 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,3,3',4,4'-PentaCB | 105 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,3,3',4,6-PentaCB | 110 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,3',4,4',5-PentaCB | 118 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,4,5,5'-HexaCB | 146 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,3,3',4,4',5-HexaCB | 156 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| 2,2',3,4,5,5',6-HeptaCB | 187 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| DecaCB | 209 | 0.2 | 1 | 5 | 50 | 400 | 2000 |
| Internal | | | | | | | |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 25 | 25 | 25 | 25 | 25 | 25 |
| Recovery | | | | | | | |
| 2,3',4',5-TetraCB | 70 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3,3',5,5'-PentaCB | 111 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',4,4',6,6'-HexaCB | 155 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3',4,4',5,5'-HexaCB | 167 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 25 | 25 | 25 | 25 | 25 | 25 |

WHO "Dioxin-Like" and "Non-Dioxin-Like" (Marker/Indicator) PCB Mixtures

| Catalog No. | Compound | | Amount |
|--------------------|---|-------|--------------------|
| NEW EC-5532 | PCB Exhibit Internal Standard | | 1.2 mL in nonane |
| | Labeled | IUPAC | (ng/mL) |
| | 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | 52 | 100 |
| | 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 100 |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 |
| | 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 100 |
| | 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 100 |
| | 2,2',3,3',4,4'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 100 |
| | 2,3,3',4,4',5-HexaCB (¹³ C ₁₂ , 99%) | 156 | 100 |
| | 2,3,3',4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 157 | 100 |
| | 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 100 |
| | 2,2',3,4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 180 | 100 |
| | DecaCB (¹³ C ₁₂ , 99%) | 209 | 100 |
| NEW EC-5533 | PCB Exhibit Recovery Solution | | 1.2 mL in nonane |
| | Labeled | IUPAC | (ng/mL) |
| | 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 5000 |
| | 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 5000 |
| | 2,2',4,4',6,6'-HexaCB (¹³ C ₁₂ , 99%) | 155 | 5000 |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 5000 |
| | 2,3,3',4,4',5,5'-HeptaCB (¹³ C ₁₂ , 99%) | 189 | 5000 |
| NEW EC-5534 | PCB Exhibit Native PAR Solution | | 1.2 mL in isoctane |
| | Unlabeled | IUPAC | (ng/mL) |
| | 2,2',5,5'-TetraCB | 52 | 200 |
| | 2,3',4,4'-TetraCB | 66 | 200 |
| | 2,4,4',5-TetraCB | 74 | 200 |
| | 2,2',3,4,5'-PentaCB | 87 | 200 |
| | 2,2',4,4',5-PentaCB | 99 | 200 |
| | 2,2',4,5,5'-PentaCB | 101 | 200 |
| | 2,3,3',4,4'-PentaCB | 105 | 200 |
| | 2,3,3',4',6-PentaCB | 110 | 200 |
| | 2,3',4,4',5-PentaCB | 118 | 200 |
| | 2,2',3,4,4',5'-HexaCB | 138 | 200 |
| | 2,2',3,4',5,5'-HexaCB | 146 | 200 |
| | 2,2',4,4',5,5'-HexaCB | 153 | 200 |
| | 2,3,3',4,4',5-HexaCB | 156 | 200 |
| | 2,2',3,3',4,4',5-HeptaCB | 170 | 200 |
| | 2,2',3,4,4',5,5'-HeptaCB | 180 | 200 |
| | 2,2',3,4',5,5',6-HeptaCB | 187 | 200 |
| | DecaCB | 209 | 200 |

Isotope-Labeled PCB Standard Mixtures

| Catalog No. | Compound | | Amount |
|-----------------|---|-------|---|
| EC-5181 | PCB Syringe Spike | | 1.2 mL in nonane |
| EC-5181-10X-1.2 | PCB Syringe Spike | | 1.2 mL in nonane |
| | | | |
| | Labeled | IUPAC | EC-5181 (ng/mL) EC-5181-10X-1.2 (ng/mL) |
| | 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 100 1000 |
| | 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 100 1000 |
| | 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 100 1000 |
| | | | |
| EC-5163 | PCB Mixture (PCB-70/111/138/1 70) | | 1.2 mL in nonane |
| | | | |
| | Labeled | IUPAC | (ng/mL) |
| | 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 1000 |
| | 2,3,3',5,5'-PentaCB (¹³ C ₁₂ , 99%) | 111 | 1000 |
| | 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | 138 | 1000 |
| | 2,2',3,3',4,4',5-HeptaCB (¹³ C ₁₂ , 99%) | 170 | 1000 |
| | | | |
| EC-5350 | POPs Pesticides HRMS (PCB) Syringe Spike | | 1.2 mL in nonane |
| EC-5350-L | POPs Pesticides HRMS (PCB) Syringe Spike | | 1.2 mL in nonane |
| | | | |
| | Labeled | IUPAC | EC-5350 (ng/mL) EC-5350-L (ng/mL) |
| | 4,4'-DiCB (¹³ C ₁₂ , 99%) | 15 | 100 1000 |
| | 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) | 70 | 100 1000 |
| | | | |
| NEW EC-5523 | PCB Recovery Standard (PCB-9, 118, 188) | | 1.2 mL in nonane |
| | | | |
| | Labeled | IUPAC | (ng/mL) |
| | 2,5-DiCB (¹³ C ₁₂ , 99%) | 9 | 12,500 |
| | 2,3',4,4',5-PentaCB (¹³ C ₁₂ , 99%) | 118 | 12,500 |
| | 2,2',3,4',5,6,6'-HeptaCB (¹³ C ₁₂ , 99%) | 188 | 12,500 |
| | | | |
| NEW ES-5474 | CDC PCB Recovery Standard for OH-PAHs | | 1 mL in toluene |
| | | | |
| | Labeled | IUPAC | (ng/mL) |
| | 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 200 |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 200 |
| | 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 200 |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 200 |
| | | | |
| EC-4060 | PCB Mixture | | 1.2 mL in nonane |
| | | | |
| | Labeled | IUPAC | (ng/mL) |
| | 3,3',4,4'-TetraCB (¹³ C ₁₂ , 99%) | 77 | 10,000 |
| | 2,2',4,5,5'-PentaCB (¹³ C ₁₂ , 99%) | 101 | 10,000 |
| | 2,2',3,4,5,5'-HexaCB (¹³ C ₁₂ , 99%) | 141 | 10,000 |
| | 2,2',3,3',5,5',6-HeptaCB (¹³ C ₁₂ , 99%) | 178 | 10,000 |

Unlabeled PCB Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|--------------------|
| EC-5434 | Fully Resolved Native Mono-Deca PCB Mixture | 1.2 mL in isoctane |

| Unlabeled | IUPAC | (ng/mL) |
|-------------------|-------|---------|
| 2-MonoCB | 1 | 2000 |
| 4-MonoCB | 3 | 2000 |
| 2,4'-DiCB | 8 | 2000 |
| 2,5-DiCB | 9 | 2000 |
| 2,6-DiCB | 10 | 2000 |
| 3,4-DiCB | 12 | 2000 |
| 4,4'-DiCB | 15 | 2000 |
| 2,2',5-TriCB | 18 | 1000 |
| 2,2',6-TriCB | 19 | 1000 |
| 2',3,4-TriCB | 33 | 1000 |
| 3,3',4-TriCB | 35 | 1000 |
| 3,4,4'-TriCB | 37 | 1000 |
| 3,4,5-TriCB | 38 | 1000 |
| 2,2',3,5'-TetraCB | 44 | 1000 |
| 2,2',5,5'-TetraCB | 52 | 1000 |
| 2,2',6,6'-TetraCB | 54 | 1000 |
| 2,3,3',5-TetraCB | 57 | 1000 |
| 2,4,4',5-TetraCB | 74 | 1000 |
| 3,3',4,4'-TetraCB | 77 | 1000 |
| 3,3',4,5-TetraCB | 78 | 1000 |
| 3,3',4,5'-TetraCB | 79 | 1000 |
| 3,4,4'-TetraCB | 81 | 1000 |

| Unlabeled | IUPAC | (ng/mL) |
|------------------------------|-------|---------|
| 2,2',4,4',5-PentaCB | 99 | 1000 |
| 2,2',4,6,6'-PentaCB | 104 | 1000 |
| 2,3,4,4',5-PentaCB | 114 | 1000 |
| 2,3',4,4',5-PentaCB | 118 | 1000 |
| 2',3,4,4',5-PentaCB | 123 | 1000 |
| 3,3',4,4',5-PentaCB | 126 | 1000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 1000 |
| 2,2',4,4',6,6'-HexaCB | 155 | 1000 |
| 2,3,3',4,4',5-HexaCB | 156 | 1000 |
| 2,3,3',4,4',5'-HexaCB | 157 | 1000 |
| 2,3,3',4,5,5'-HexaCB | 162 | 1000 |
| 2,3',4,4',5,5'-HexaCB | 167 | 1000 |
| 3,3',4,4',5,5'-HexaCB | 169 | 1000 |
| 2,2',3,4',5,6,6'-HeptaCB | 188 | 1000 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 1000 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 1000 |
| 2,2',3,3',4,4',5,6-OctaCB | 195 | 1000 |
| 2,2',3,3',5,5',6,6'-OctaCB | 202 | 1000 |
| 2,3,3',4,4',5,5',6-OctaCB | 205 | 1000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 1000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 1000 |
| DecaCB | 209 | 1000 |

| | | |
|---------|-----------------|------------------|
| EC-4133 | DSJ PCB Mixture | 1 mL in isoctane |
|---------|-----------------|------------------|

| Unlabeled | IUPAC | (ng/mL) |
|-----------------------|-------|---------|
| 2,2',4-TriCB | 17 | 500 |
| 2,2',5-TriCB | 18 | 2000 |
| 2,4,4'-TriCB | 28 | 2000 |
| 2,4',5-TriCB | 31 | 1500 |
| 2',3,4-TriCB | 33 | 2000 |
| 2,2',3,5'-TetraCB | 44 | 2000 |
| 2,2',4,5'-TetraCB | 49 | 2000 |
| 2,2',5,5'-TetraCB | 52 | 2000 |
| 2,3',4',5-TetraCB | 70 | 2000 |
| 2,4,4',5-TetraCB | 74 | 2000 |
| 2,2',3,3',4-PentaCB | 82 | 500 |
| 2,2',3,4,5'-PentaCB | 87 | 2000 |
| 2,2',3,5',6-PentaCB | 95 | 1000 |
| 2,2',4,4',5-PentaCB | 99 | 2000 |
| 2,2',4,5,5'-PentaCB | 101 | 2000 |
| 2,3,3',4,4'-PentaCB | 105 | 500 |
| 2,3,3',4',6-PentaCB | 110 | 2000 |
| 2,3',4,4',5-PentaCB | 118 | 2000 |
| 2,2',3,3',4,4'-HexaCB | 128 | 2000 |
| 2,2',3,3',4,6'-HexaCB | 132 | 1000 |
| 2,2',3,4,4',5'-HexaCB | 138 | 2000 |

| Unlabeled | IUPAC | (ng/mL) |
|------------------------------|-------|---------|
| 2,2',3,4',5',6-HexaCB | 149 | 2000 |
| 2,2',3,5,5',6-HexaCB | 151 | 2000 |
| 2,2',4,4',5,5'-HexaCB | 153 | 2000 |
| 2,3,3',4,4',5-HexaCB | 156 | 2000 |
| 2,3,3',4,4',6-HexaCB | 158 | 500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 2000 |
| 2,2',3,3',4,4',5-HeptaCB | 170 | 2000 |
| 2,2',3,3',4,4',6-HeptaCB | 171 | 2000 |
| 2,2',3,3',4',5,6-HeptaCB | 177 | 2000 |
| 2,2',3,4,4',5,5'-HeptaCB | 180 | 2000 |
| 2,2',3,4,4',5,6-HeptaCB | 183 | 2000 |
| 2,2',3,4',5,5',6-HeptaCB | 187 | 2000 |
| 2,3,3',4,4',5,6-HeptaCB | 191 | 2000 |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 2000 |
| 2,2',3,3',4,4',5,6-OctaCB | 195 | 2000 |
| 2,2',3,3',4,5,5',6-OctaCB | 201 | 1500 |
| 2,3,3',4,4',5,5',6-OctaCB | 205 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 2000 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 2000 |
| DecaCB | 209 | 2000 |

Unlabeled PCB Standard Mixtures

| Catalog No. | Compound | | Amount |
|---------------------|---|-------|--------------------|
| EC-5492 | PCB Recovery Standard | | 1 mL in toluene |
| | Unlabeled | IUPAC | (ng/mL) |
| | 2,4,4'-TriCB | 28 | 200 |
| | 2,3,3',4,4'-PentaCB | 105 | 200 |
| | 2,2',4,4',5,5'-HexaCB | 153 | 200 |
| | 2,3',4,4',5,5'-HexaCB | 167 | 200 |
| <hr/> | | | |
| NEW EC-5504 | PCB Recovery Standard | | 1.2 mL in isoctane |
| | Unlabeled | IUPAC | (ng/mL) |
| | 2,3',4',5-TetraCB | 70 | 100 |
| | 2,3,3',5,5'-PentaCB | 111 | 100 |
| | 2,2',3,3',4,4',5-HeptaCB | 170 | 100 |
| <hr/> | | | |
| EC-7438 | PCB Mixture | | 1.2 mL in isoctane |
| | Unlabeled | IUPAC | (ng/mL) |
| | 3,3',4,4'-TetraCB | 77 | 10,000 |
| | 2,2',4,5,5'-PentaCB | 101 | 10,000 |
| | 2,2',3,4,5,5'-HexaCB | 141 | 10,000 |
| | 2,2',3,3',5,5',6-HeptaCB | 178 | 10,000 |
| <hr/> | | | |
| NEW ULM-5370 | EN-1948-4 PCB Sampling Standard | | Inquire |
| | Unlabeled | IUPAC | (ng/mL) |
| | 2,3,4,4'-TetraCB | 60 | 100 |
| | 3,3',4,5,5'-PentaCB | 127 | 100 |
| | 2,3,3',4,5,5'-HexaCB | 159 | 100 |
| <hr/> | | | |
| NEW EC-5460 | UNEP OC Pesticide ECD Internal Standard Mixture | | 1.2 mL in isoctane |
| | Unlabeled | IUPAC | (ng/mL) |
| | 2,3,3',5,6-PentaCB | 112 | 5000 |
| | 2,2',4,4',6,6'-HexaCB | 155 | 5000 |
| | 2,2',3,3',4,5,5',6-OctaCB | 198 | 5000 |

PCB Window Defining Mixture

| Catalog No. | Compound | Amount |
|---|--|------------------|
| EC-1430 | PCB Window Defining Mixture (for use with DB-5 type GC-MS columns) | 5 mL in isoctane |
| Unlabeled | | |
| Biphenyl | IUPAC | (ng/mL) |
| 2-MonoCB | 0 | 2500 |
| 4-MonoCB | 1 | 2500 |
| 2,6-DiCB | 3 | 2500 |
| 2,6-DiCB | 10 | 2500 |
| 4,4'-DiCB | 15 | 2500 |
| <i>Note: #30 is second tri eluter</i> | | |
| 2,4,6-TriCB | 30 | 2500 |
| 3,4,4'-TriCB | 37 | 2500 |
| 2,2',6,6'-TetraCB | 54 | 2500 |
| 3,3',4,4'-TetraCB | 77 | 2500 |
| 2,2',4,6,6'-PentaCB | 104 | 2500 |
| 3,3',4,4',5-PentaCB | 126 | 2500 |
| 2,2',4,4',6,6'-HexaCB | 155 | 2500 |
| 3,3',4,4',5,5'-HexaCB | 169 | 2500 |
| 2,2',3,4',5,6,6'-HeptaCB | 188 | 2500 |
| 2,3,3',4,4',5,5'-HeptaCB | 189 | 2500 |
| 2,2',3,3',5,5',6,6'-OctaCB | 202 | 2500 |
| <i>Note: #194 is second-to-last octa eluter</i> | | |
| 2,2',3,3',4,4',5,5'-OctaCB | 194 | 2500 |
| 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 2500 |
| 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 2500 |
| DecaCB | 209 | 2500 |

Isotope-Labeled Mixed Bromo/Chlorobiphenyl Standards

| Catalog No. | Compound | Concentration | Amount |
|-------------|--|-------------------------------|--------|
| ECB-5269 | 3,4-Dichloro-3',4',5'-triBB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5270 | 3,4-Dibromo-3',4'-diCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5271 | 3,4-Dibromo-3',4',5'-triCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5291 | 4'-Bromo-3,3',4,5-tetraCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5292 | 4'-Bromo-2,3',4,5-tetraCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5293 | 4'-Bromo-2,3,3',4-tetraCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5294 | 4'-Bromo-2,3,3',4,5-pentaCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |
| ECB-5339 | 4'-Bromo-3,3',4,5,5'-pentaCB ($^{13}\text{C}_{12}$, 99%) | 40 $\mu\text{g/mL}$ in nonane | 3 mL |

Unlabeled Mixed Bromo/Chlorobiphenyl Standards

| Catalog No. | Compound | Concentration | Amount |
|--------------|---|----------------------------------|--------|
| PCBB-5272-CS | 3,4-Dichloro-3',4',5'-triBB (Certified Standard) | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5273 | 3,4-Dibromo-3',4'-diCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5274 | 3,4-Dibromo-3',4',5'-triCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5295 | 4'-Bromo-3,3',4,5-tetraCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5296 | 4'-Bromo-2,3',4,5-tetraCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5297 | 4'-Bromo-2,3,3',4-tetraCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5298 | 4'-Bromo-2,3,3',4,5-pentaCB | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| PCBB-5340-CS | 4'-Bromo-3,3',4,5,5'-pentaCB (Certified Standard) | 100 $\mu\text{g/mL}$ in isoctane | 1.2 mL |

Mixed Bromo/Chlorobiphenyl Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------|-------------------------------------|-----------------------------|
| ECB-5390 | PXB Calibration Solutions [CS1-CS5] | Set of 5 × 0.2 mL in nonane |
| ECB-5390-CS1 | PXB Calibration Solution [CS1] | 0.2 mL in nonane |
| ECB-5390-CS2 | PXB Calibration Solution [CS2] | 0.2 mL in nonane |
| ECB-5390-CS3 | PXB Calibration Solution [CS3] | 0.2 mL in nonane |
| ECB-5390-CS4 | PXB Calibration Solution [CS4] | 0.2 mL in nonane |
| ECB-5390-CS5 | PXB Calibration Solution [CS5] | 0.2 mL in nonane |

| All concentrations are in ng/mL (ppb) | | | | | |
|---|-----|-----|-----|-----|------|
| Unlabeled | CS1 | CS2 | CS3 | CS4 | CS5 |
| 4'-Bromo-3,3',4,5-tetraCB | 2 | 10 | 50 | 200 | 200 |
| 4'-Bromo-2,3',4,5-tetraCB | 2 | 10 | 50 | 200 | 200 |
| 4'-Bromo-2,3,3',4-tetraCB | 2 | 10 | 50 | 200 | 200 |
| 4'-Bromo-2,3,3',4,5-pentaCB | 2 | 10 | 50 | 200 | 200 |
| 4'-Bromo-3,3',4,5,5'-pentaCB | 2 | 10 | 50 | 200 | 1000 |
| 3,4-Dichloro-3',4',5'-triBB | 4 | 20 | 100 | 400 | 2000 |
| Labeled | | | | | |
| 4'-Bromo-3,3',4,5-tetraCB (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 4'-Bromo-2,3',4,5-tetraCB (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 4'-Bromo-2,3,3',4-tetraCB (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 4'-Bromo-2,3,3',4,5-pentaCB (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 4'-Bromo-3,3',4,5,5'-pentaCB (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |
| 3,4-Dichloro-3',4',5'-triBB (¹³ C ₁₂ , 99%) | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,4,5,5'-HexaCDE (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 |

| | | |
|----------|-------------------|------------------|
| ECB-5389 | PXB Cleanup Spike | 1.2 mL in nonane |
|----------|-------------------|------------------|

| Labeled | (ng/mL) |
|---|---------|
| 4'-Bromo-3,3',4,5-tetraCB (¹³ C ₁₂ , 99%) | 1000 |
| 4'-Bromo-2,3',4,5-tetraCB (¹³ C ₁₂ , 99%) | 1000 |
| 4'-Bromo-2,3,3',4-tetraCB (¹³ C ₁₂ , 99%) | 1000 |
| 4'-Bromo-2,3,3',4,5-pentaCB (¹³ C ₁₂ , 99%) | 1000 |
| 4'-Bromo-3,3',4,5,5'-pentaCB (¹³ C ₁₂ , 99%) | 1000 |
| 3,4-Dichloro-3',4',5'-triBB (¹³ C ₁₂ , 99%) | 2000 |

| | | |
|--|----------------------|------------------|
| EO-5388 | PXB Syringe Standard | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| 2,2',3,4,5,5'-HexaCDE (¹³ C ₁₂ , 99%) | 1000 | |

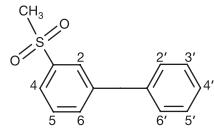
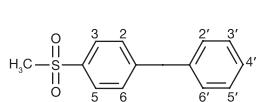
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|------------------------------|-------------------------|------------------|
| ECB-5387 | PXB Native PAR Solution | 0.5 mL in nonane |
| Unlabeled (ng/mL) | | |
| 4'-Bromo-3,3',4,5-tetraCB | 1000 | |
| 4'-Bromo-2,3',4,5-tetraCB | 1000 | |
| 4'-Bromo-2,3,3',4-tetraCB | 1000 | |
| 4'-Bromo-2,3,3',4,5-pentaCB | 1000 | |
| 4'-Bromo-3,3',4,5,5'-pentaCB | 1000 | |
| 3,4-Dichloro-3',4',5'-triBB | 2000 | |

PCB Metabolite Standards

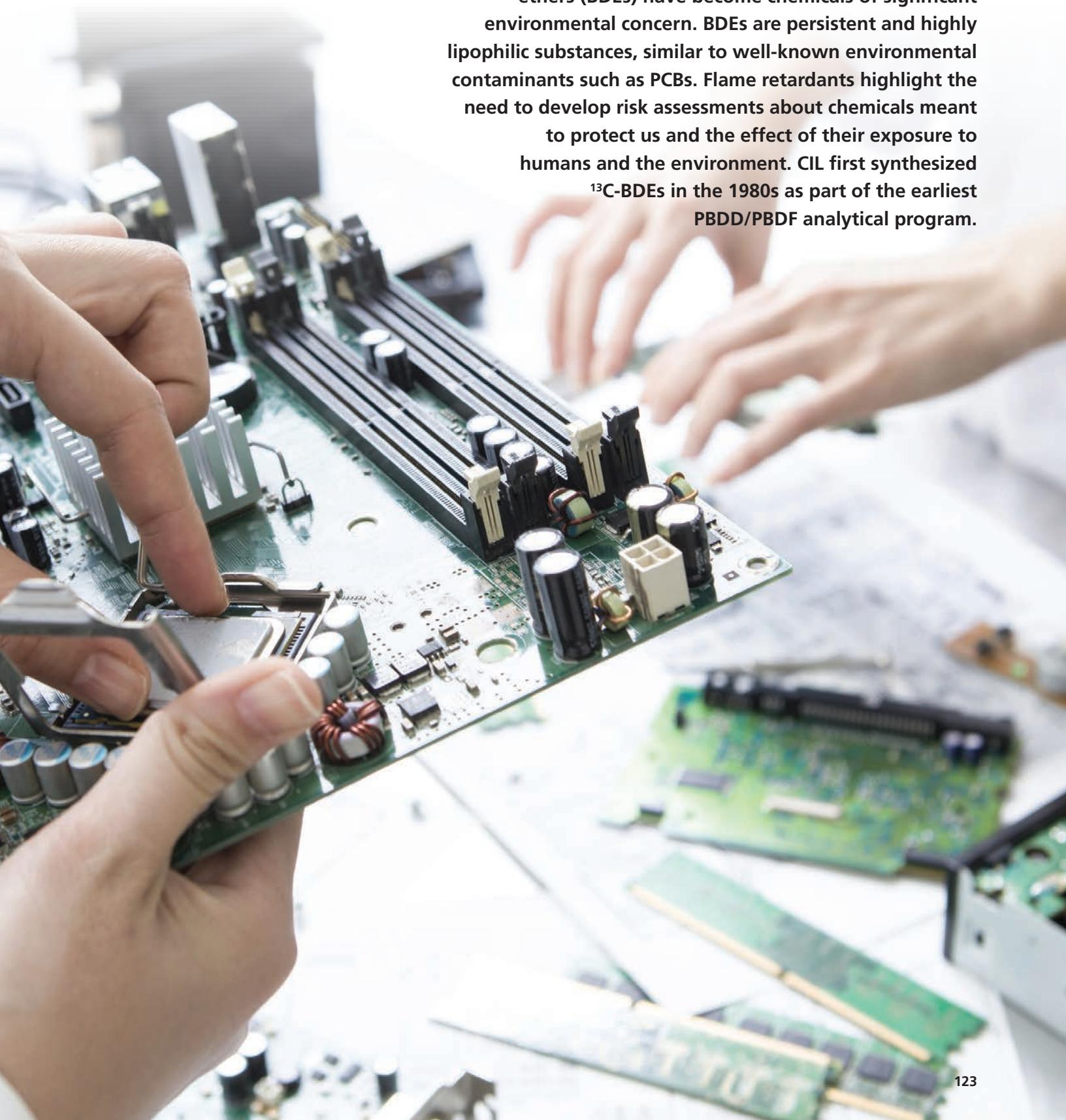
| Catalog No. | Compound | Concentration | Amount |
|---------------------------|---|--------------------|--------|
| OH-PCB | | | |
| OHCB-5114-1.2 | 4'-OH-3,3',4,5'-tetraCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| OHCB-5115-1.2 | 4-OH-2,3,3',4',5-pentaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| OHCB-5117-1.2 | 4-OH-2,2',3,4',5,5'-hexaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| OHCB-5118-1.2 | 3'-OH-2,2',3,4,4',5'-hexaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| OHCB-5124-1.2 | 4'-OH-2,2',3,3',4,5,5'-heptaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| MeO-PCB | | | |
| <i>NEW</i> MEOCB-5485-1.2 | 4-Methoxy-2,3,3',4',5-pentaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOCB-5486-1.2 | 4-Methoxy-2,2',3,4',5,5'-hexaCB ($^{13}\text{C}_{12}$, 99%) | 50 µg/mL in nonane | 1.2 mL |
| MEOCB-5109-1.2 | 4-Methoxy-2,3,3',4',5-pentaCB (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| MEOCB-5111-1.2 | 4-Methoxy-2,2',3,4',5,5'-hexaCB (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| MEOCB-5135-1.2 | 4-Methoxy-2,2',3,4',5,5',6-heptaCB (unlabeled) | 50 µg/mL in nonane | 1.2 mL |

Unlabeled Methyl Sulfone PCB Standards

| Catalog No. | Compound | Concentration | Amount |
|-------------|---|------------------------|--------|
| MSCB-4027 | 3-MeSO ₂ -4-Me-2',3',4',5,5'-pentaCB (Internal Standard) | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4007 | 3-MeSO ₂ -2,2',4',5-tetraCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4008 | 4-MeSO ₂ -2,2',4',5-tetraCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4009 | 3-MeSO ₂ -2,2',4',5,5'-pentaCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4010 | 4-MeSO ₂ -2,2',4',5,5'-pentaCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4012 | 4-MeSO ₂ -2,3,3',4',6-pentaCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4013 | 3-MeSO ₂ -2,2',4',5,5',6-hexaCB | 40 ± 4 µg/mL in nonane | 1.2 mL |
| MSCB-4015 | 3-MeSO ₂ -DDE | 40 ± 4 µg/mL in nonane | 1.2 mL |

3-MeSO₂-PCB4-MeSO₂-PCB

Generally used as flame retardants, brominated diphenyl ethers (BDEs) have become chemicals of significant environmental concern. BDEs are persistent and highly lipophilic substances, similar to well-known environmental contaminants such as PCBs. Flame retardants highlight the need to develop risk assessments about chemicals meant to protect us and the effect of their exposure to humans and the environment. CIL first synthesized ^{13}C -BDEs in the 1980s as part of the earliest PBDD/PBDF analytical program.



Isotope-Labeled and Unlabeled Brominated Diphenyl Ether (BDE) Standards

Generally used as flame retardants, brominated diphenyl ethers (BDEs) have become chemicals of significant environmental concern. BDEs may be generally considered as persistent and highly lipophilic substances, similar to well known environmental contaminants such as polychlorinated biphenyls (PCBs).

In 1996, CIL first introduced chlorinated and brominated diphenyl ether analytical standards. Since then, additional studies in Europe, Canada, Japan, and the United States have significantly increased interest in these products and enabled CIL to extend its flame-retardant offerings.

US EPA Method 1614 Standard Mixtures

The US EPA has developed a standardized test method for analysis of BDEs in multiple matrices by high-resolution GC/MS. CIL worked closely with the EPA and their contracting laboratories to develop standard mixtures specifically for use in Method 1614.

RoHS BDE Standard Mixtures

The Directive on the Restriction of the Use of Certain Hazardous Substances in Electrical and Electronic Equipment (2002/95/EC), commonly referred to as the Restriction of Hazardous Substances Directive (RoHS) was adopted in February 2003 in Europe and took effect in July 2006. This directive restricts the use of several types of hazardous materials in the manufacture of various types of electronic and electrical equipment. CIL worked with laboratories in the European Union to develop standards for analysis of brominated diphenyl ethers under the RoHS guidelines.

Isotope-Labeled and Unlabeled BDE Metabolites

Researchers have suggested that BDE body burdens are not completely represented by measurements of BDEs in tissue or milk. Analytical data indicates that the liver hydrolyzes BDEs in its attempts to expel them. BDE toxicity is still being established, but there is a likelihood that BDE metabolites have similar or greater toxicity than the parent BDEs. CIL has been producing both unlabeled and ¹³C₁₂ methoxy- and hydroxy-BDEs. These items represent some of the BDE metabolites available from CIL; please contact CIL or your CIL representative for more information on these compounds.

Brominated Flame-Retardant Standards and Standard Mixtures

Following regulatory restrictions and voluntary removal of BDE products from many flame-retardant applications, new brominated compounds have been developed as replacements for BDEs. As monitoring of these possible environmental pollutants increased, CIL developed several labeled and unlabeled standards to allow accurate analysis of these new-use brominated flame retardant (BFR) compounds. CIL further developed a series of standard mixtures, which include a combination of BDEs and other BFRs in comprehensive mixtures.

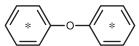
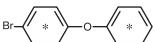
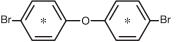
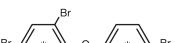
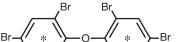
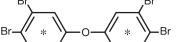
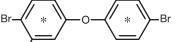
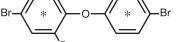
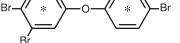
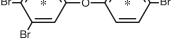
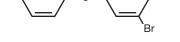
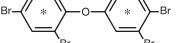
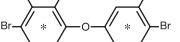
Dechloranes

Dechloranes are highly chlorinated norbornene compounds, synthesized by the Diels-Alder reaction of hexachlorocyclopenta-diene with various substrates. Dechloranes first found commercial application as flame-retardant compounds in the 1960s. While Dechlorane Plus has been the subject of significant environmental research in recent years, related compounds – Dechlorane 602 (Dec 602), Dec 603 and Dec 604 – have been identified in air, fish and sediment samples. Although the research thus far is limited, indications suggest that Dec 602 may be found in fish and sediment samples at levels similar to or even higher than Dechlorane Plus.

Phosphorus Flame Retardants

Since the 1940s, trialkyl, triaryl and halogenated trialkyl/triaryl phosphate ester compounds have been used as commercial additives in flame retardants, plasticizers, hydraulic fluids, solvents, extraction agents, antifoam agents, adhesives and coatings for electronic devices. In recent years, the demand for phosphate ester compounds for flame-retardant applications has increased as many brominated flame retardants have been removed from the market. New research initiatives continue to broaden the understanding of the effects of these compounds in the environment, though analytical challenges remain. The use of isotopically labeled internal standards has been shown to help compensate for matrix effects, and to generally strengthen data by use of isotope dilution mass spectrometry.

Isotope-Labeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|--------------|---|-----|--|------------------------|--------|
| CLM-1587-1.2 | Diphenyl ether ($^{13}\text{C}_{12}$, 99%) | 0 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4999 | 4-Monobromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 3 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5001 | 4,4'-Dibromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 15 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5002 | 2,4,4'-Tribromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 28 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4982 | 2,2',4,4'-Tetrabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 47 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-1439 | 3,3',4,4'-Tetrabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 77 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4983 | 2,2',4,4',5-Pentabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 99 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4993 | 2,2',4,4',6-Pentabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 100 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5034 | 2,3',4,4',5-Pentabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 118 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4930 | 3,3',4,4',5-Pentabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 126 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5035 | 2,2',3,4,4',5'-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 138 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5165 | 2,2',3,4,4',6-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 139 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4984 | 2,2',4,4',5,5'-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 153 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5161 | 2,2',4,4',5,6'-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 154 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5413 | 2,2',4,4',6,6'-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 155 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-4985 | 2,2',3,4,4',5',6-Heptabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 183 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5376 | 2,3,3',4,4',5,6-Heptabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 190 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual ^{13}C -BDE standards as follows: 2-3 units = 20% discount; 4 or more units = 33% discount

(continued)

Isotope-Labeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|-------------|---|-----|-----------|------------------------|--------|
| EO-5337 | 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether (¹³ C ₁₂ , 99%) | 197 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5377 | 2,2',3,4,4',5,5',6-Octabromodiphenyl ether (¹³ C ₁₂ , 99%) | 203 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5355 | 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether (¹³ C ₁₂ , 99%) | 204 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5362 | 2,3,3',4,4',5,5',6-Octabromodiphenyl ether (¹³ C ₁₂ , 99%) | 205 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5240 | 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether (¹³ C ₁₂ , 99%) | 206 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5241 | 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether (¹³ C ₁₂ , 99%) | 207 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5242 | 2,2',3,3',4,5,5',6,6'-Nonabromodiphenyl ether (¹³ C ₁₂ , 99%) | 208 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| EO-5003 | Decabromodiphenyl ether (¹³ C ₁₂ , 99%) | 209 | | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual ¹³C-BDE standards as follows: 2-3 units = 20% discount; 4 or more units = 33% discount

Unlabeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|-------------|----------------------------|-----|-----------|------------------------|--------|
| BDE-1-CS | 2-Monobromodiphenyl ether | 1 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-2-CS | 3-Monobromodiphenyl ether | 2 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-3-CS | 4-Monobromodiphenyl ether | 3 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-7-CS | 2,4-Dibromodiphenyl ether | 7 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-8-CS | 2,4'-Dibromodiphenyl ether | 8 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-10-CS | 2,6-Dibromodiphenyl ether | 10 | | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual unlabeled BDE standards as follows: 5-7 units = 20% discount; 8-11 units = 33% discount; 12 or more units = 50% discount

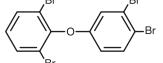
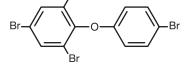
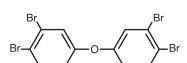
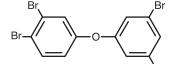
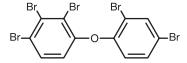
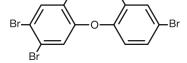
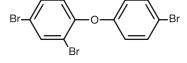
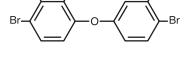
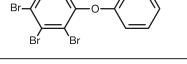
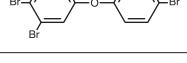
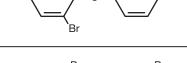
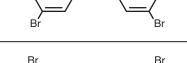
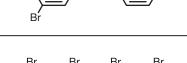
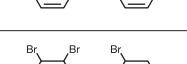
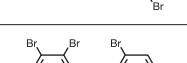
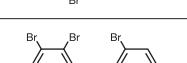
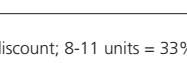
Unlabeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|-------------|------------------------------------|-----|-----------|------------------------|--------|
| BDE-11-CS | 3,3'-Dibromodiphenyl ether | 11 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-12-CS | 3,4-Dibromodiphenyl ether | 12 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-13-CS | 3,4'-Dibromodiphenyl ether | 13 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-15-CS | 4,4'-Dibromodiphenyl ether | 15 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-17-CS | 2,2',4-Tribromodiphenyl ether | 17 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-25-CS | 2,3',4-Tribromodiphenyl ether | 25 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-28-CS | 2,4,4'-Tribromodiphenyl ether | 28 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-30-CS | 2,4,6-Tribromodiphenyl ether | 30 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-32-CS | 2,4',6-Tribromodiphenyl ether | 32 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-33-CS | 2',3,4-Tribromodiphenyl ether | 33 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-35-CS | 3,3',4-Tribromodiphenyl ether | 35 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-37-CS | 3,4,4'-Tribromodiphenyl ether | 37 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-47-CS | 2,2',4,4'-Tetrabromodiphenyl ether | 47 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-49-CS | 2,2',4,5'-Tetrabromodiphenyl ether | 49 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-51-CS | 2,2',4,6'-Tetrabromodiphenyl ether | 51 | | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-66-CS | 2,3',4,4'-Tetrabromodiphenyl ether | 66 | | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual unlabeled BDE standards as follows: 5-7 units = 20% discount; 8-11 units = 33% discount; 12 or more units = 50% discount

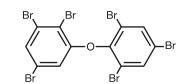
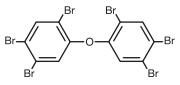
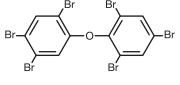
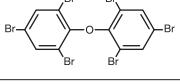
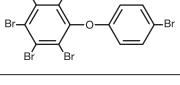
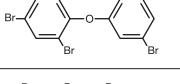
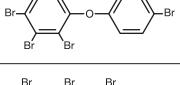
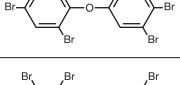
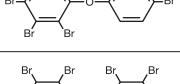
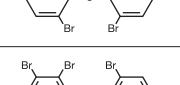
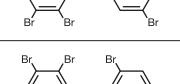
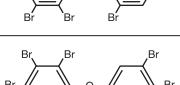
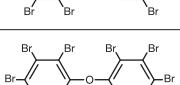
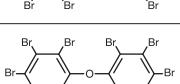
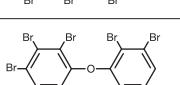
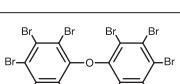
(continued)

Unlabeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|-------------|--|-----|--|------------------------|--------|
| BDE-71-CS | 2,3',4',6-Tetrabromodiphenyl ether | 71 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-75-CS | 2,4,4',6-Tetrabromodiphenyl ether | 75 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-77-CS | 3,3',4,4'-Tetrabromodiphenyl ether | 77 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-79-CS | 3,3',4,5'-Tetrabromodiphenyl ether | 79 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-85-CS | 2,2',3,4,4'-Pentabromodiphenyl ether | 85 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-99-CS | 2,2',4,4',5-Pentabromodiphenyl ether | 99 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-100-CS | 2,2',4,4',6-Pentabromodiphenyl ether | 100 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-105-CS | 2,3,3',4,4'-Pentabromodiphenyl ether | 105 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-116-CS | 2,3,4,5,6-Pentabromodiphenyl ether | 116 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-118-CS | 2,3',4,4',5-Pentabromodiphenyl ether | 118 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-119-CS | 2,3',4,4',6-Pentabromodiphenyl ether | 119 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-120-CS | 2,3',4,5,5'-Pentabromodiphenyl ether | 120 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-126-CS | 3,3',4,4',5-Pentabromodiphenyl ether | 126 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-128-CS | 2,2',3,3',4,4'-Hexabromodiphenyl ether | 128 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-138-CS | 2,2',3,4,4',5'-Hexabromodiphenyl ether | 138 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-139-CS | 2,2',3,4,4',6-Hexabromodiphenyl ether | 139 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-140-CS | 2,2',3,4,4',6'-Hexabromodiphenyl ether | 140 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual unlabeled BDE standards as follows: 5-7 units = 20% discount; 8-11 units = 33% discount; 12 or more units = 50% discount

Unlabeled Individual Brominated Diphenyl Ether (BDE) Standards

| Catalog No. | Compound | BDE | Structure | Concentration | Amount |
|-------------|---|-----|--|------------------------|--------|
| BDE-148-CS | 2,2',3,4',5,6'-Hexabromodiphenyl ether | 148 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-153-CS | 2,2',4,4',5,5'-Hexabromodiphenyl ether | 153 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-154-CS | 2,2',4,4',5,6'-Hexabromodiphenyl ether | 154 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-155-CS | 2,2',4,4',6,6'-Hexabromodiphenyl ether | 155 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-166-CS | 2,3,4,4',5,6-Hexabromodiphenyl ether | 166 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-175-CS | 2,2',3,3',4,5',6-Heptabromodiphenyl ether | 175 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-181-CS | 2,2',3,4,4',5,6-Heptabromodiphenyl ether | 181 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-183-CS | 2,2',3,4,4',5',6-Heptabromodiphenyl ether | 183 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-190-CS | 2,3,3',4,4',5,6-Heptabromodiphenyl ether | 190 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-197-CS | 2,2',3,3',4,4',6,6'-Octabromodiphenyl ether | 197 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-203-CS | 2,2',3,4,4',5,5',6-Octabromodiphenyl ether | 203 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-204-CS | 2,2',3,4,4',5,6,6'-Octabromodiphenyl ether | 204 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-205-CS | 2,3,3',4,4',5,5',6-Octabromodiphenyl ether | 205 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-206-CS | 2,2',3,3',4,4',5,5',6-Nonabromodiphenyl ether | 206 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-207-CS | 2,2',3,3',4,4',5,6,6'-Nonabromodiphenyl ether | 207 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-208-CS | 2,2',3,3',4,4,5,5',6,6'-Nonabromodiphenyl ether | 208 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |
| BDE-209-CS | Decabromodiphenyl ether | 209 |  | 50 ± 5 µg/mL in nonane | 1.2 mL |

Multiple-unit pricing is available for individual unlabeled BDE standards as follows: 5-7 units = 20% discount; 8-11 units = 33% discount; 12 or more units = 50% discount

US EPA Method 1614 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EO-5279 | EPA Method 1614 Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EO-5279-CS1 | EPA Method 1614 Calibration Solution [CS1] | 0.2 mL in nonane |
| EO-5279-CS2 | EPA Method 1614 Calibration Solution [CS2] | 0.2 mL in nonane |
| EO-5279-CS3 | EPA Method 1614 Calibration Solution [CS3] | 0.2 mL in nonane |
| EO-5279-CS4 | EPA Method 1614 Calibration Solution [CS4] | 0.2 mL in nonane |
| EO-5279-CS5 | EPA Method 1614 Calibration Solution [CS5] | 0.2 mL in nonane |

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | | |
|--|----------|------|------|------|------|--------|
| Unlabeled | Congener | CS1 | CS2 | CS3 | CS4 | CS5 |
| 2,4,4'-TriBDE | 28 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',4,4'-TetraBDE | 47 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',4,4',5-PentabDE | 99 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',4,4',6-PentabDE | 100 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 1 | 5 | 50 | 500 | 2500 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 1 | 5 | 50 | 500 | 2500 |
| DecaBDE | 209 | 10 | 50 | 500 | 5000 | 25,000 |
| Surrogate | | | | | | |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5-PentabDE (¹³ C ₁₂ , 99%) | 99 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',6-PentabDE (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 100 | 100 | 100 | 100 | 100 |
| DecaBDE (¹³ C ₁₂ , 99%) | 209 | 1000 | 1000 | 1000 | 1000 | 1000 |
| Cleanup | | | | | | |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 | 100 | 100 | 100 | 100 | 100 |
| Injection Internal | | | | | | |
| 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | PCB-52 | 100 | 100 | 100 | 100 | 100 |
| 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | PCB-138 | 100 | 100 | 100 | 100 | 100 |

| | | | |
|------------|-------------|-------------------------------|-------|
| NEW | EO-1614-KIT | EPA Method 1614 "Starter Kit" | 1 Kit |
|------------|-------------|-------------------------------|-------|

Contains one each of the following items:

| | |
|---------|---|
| EO-5279 | Method 1614 Calibration Solutions CS1-CS5 |
| EO-5277 | Method 1614 Labeled Surrogate Stock Solution |
| EO-5276 | Method 1614 Labeled Cleanup Stock Solution |
| EO-5275 | Method 1614 Labeled Injection Internal Stock Solution |
| EO-5278 | Method 1614 Native PAR Stock Solution |

US EPA Method 1614 Standard Mixtures

| Catalog No. | Compound | Amount | |
|-------------|--|------------------|--------|
| EO-5277 | EPA Method 1614 Labeled Surrogate Stock Solution | 1.2 mL in nonane | |
| | Labeled | | |
| | 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | Congener (ng/mL) | |
| | 28 | 1000 | |
| | 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 1000 |
| | 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 1000 |
| | 2,2',4,4',6-PentaBDE (¹³ C ₁₂ , 99%) | 100 | 1000 |
| | 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 1000 |
| | 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 1000 |
| | 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 1000 |
| | DecaBDE (¹³ C ₁₂ , 99%) | 209 | 10,000 |
| EO-5276 | EPA Method 1614 Labeled Cleanup Stock Solution | 1.2 mL in nonane | |
| | Labeled | | |
| | 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | Congener (ng/mL) | |
| | 139 | 1000 | |
| EO-5275 | EPA Method 1614 Labeled Injection Internal Stock Solution | 1.2 mL in nonane | |
| | Labeled | | |
| | 2,2',5,5'-TetraCB (¹³ C ₁₂ , 99%) | Congener (ng/mL) | |
| | PCB-52 | 5000 | |
| | 2,2',3,4,4',5'-HexaCB (¹³ C ₁₂ , 99%) | PCB-138 | 5000 |
| EO-5278 | EPA Method 1614 Native PAR Stock Solution | 1.2 mL in nonane | |
| | Unlabeled | | |
| | 2,4,4'-TriBDE | Congener (ng/mL) | |
| | 28 | 1000 | |
| | 2,2',4,4'-TetraBDE | 47 | 1000 |
| | 2,2',4,4',5-PentaBDE | 99 | 1000 |
| | 2,2',4,4',6-PentaBDE | 100 | 1000 |
| | 2,2',4,4',5,5'-HexaBDE | 153 | 1000 |
| | 2,2',4,4',5,6'-HexaBDE | 154 | 1000 |
| | 2,2',3,4,4',5',6-HeptaBDE | 183 | 1000 |
| | DecaBDE | 209 | 10,000 |

RoHS Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EO-5402 | RoHS PBDE Calibration Solutions [CS1-CS5] | Set of 5 x 0.2 mL in nonane |
| EO-5402-CS1 | RoHS PBDE Calibration Solution [CS1] | 0.2 mL in nonane |
| EO-5402-CS2 | RoHS PBDE Calibration Solution [CS2] | 0.2 mL in nonane |
| EO-5402-CS3 | RoHS PBDE Calibration Solution [CS3] | 0.2 mL in nonane |
| EO-5402-CS4 | RoHS PBDE Calibration Solution [CS4] | 0.2 mL in nonane |
| EO-5402-CS5 | RoHS PBDE Calibration Solution [CS5] | 0.2 mL in nonane |

| Unlabeled | BDE | <i>All concentrations are in ng/mL (ppb)</i> | | | | |
|--|-----|--|-----|-----|-----|------|
| | | CS1 | CS2 | CS3 | CS4 | CS5 |
| 4-MonoBDE | 3 | 1 | 5 | 20 | 100 | 500 |
| 2,4-DiBDE | 7 | 1 | 5 | 20 | 100 | 500 |
| 4,4'-DiBDE | 15 | 1 | 5 | 20 | 100 | 500 |
| 2,2',4-TriBDE | 17 | 1 | 5 | 20 | 100 | 500 |
| 2,4,4'-TriBDE | 28 | 1 | 5 | 20 | 100 | 500 |
| 2,2',4,4'-TetraBDE | 47 | 1 | 5 | 20 | 100 | 500 |
| 2,2',4,5'-TetraBDE | 49 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4'-TetraBDE | 66 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,6-TetraBDE | 71 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4'-TetraBDE | 77 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,4,4'-PentabDE | 85 | 1 | 5 | 20 | 100 | 500 |
| 2,2',4,4',5-PentabDE | 99 | 1 | 5 | 20 | 100 | 500 |
| 2,2',4,4',6-PentabDE | 100 | 1 | 5 | 20 | 100 | 500 |
| 2,3',4,4',6-PentaBDE | 119 | 1 | 5 | 20 | 100 | 500 |
| 3,3',4,4',5-PentaBDE | 126 | 1 | 5 | 20 | 100 | 500 |
| 2,2',3,4,4',5'-HexaBDE | 138 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',4,4',6,6'-HexaBDE | 155 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,4,4',5,6-HexaBDE | 166 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',3,4,4',5,6-HeptaBDE | 181 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4',5,6-HeptaBDE | 190 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',3,4,4',5,5',6-OctaBDE | 203 | 2 | 10 | 40 | 200 | 1000 |
| 2,3,3',4,4',5,5',6-OctaBDE | 205 | 2 | 10 | 40 | 200 | 1000 |
| 2,2',3,3',4,4',5,5',6-NonaBDE | 206 | 5 | 25 | 100 | 500 | 2500 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE | 207 | 5 | 25 | 100 | 500 | 2500 |
| DecaBDE | 209 | 5 | 25 | 100 | 500 | 2500 |
| Cleanup | | | | | | |
| 4-MonoBDE (¹³ C ₁₂ , 99%) | 3 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5-PentabDE (¹³ C ₁₂ , 99%) | 99 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 200 | 200 | 200 | 200 | 200 |
| 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,4,4',5,6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,4,4',5,6,6'-OctaBDE (¹³ C ₁₂ , 99%) | 204 | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 207 | 500 | 500 | 500 | 500 | 500 |
| DecaBDE (¹³ C ₁₂ , 99%) | 209 | 500 | 500 | 500 | 500 | 500 |
| Syringe | | | | | | |
| 2,2',3,4,4',5'-HexaBDE (¹³ C ₁₂ , 99%) | 138 | 200 | 200 | 200 | 200 | 200 |

RoHS Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------------------|---|------------------|
| EO-5403 | RoHS PBDE Cleanup Spike | 1.2 mL in nonane |
| NEW EO-5403-10X-1.2 | RoHS PBDE Cleanup 10X Spike ($^{13}\text{C}_{12}$, 99%) | 1.2 mL in nonane |

| Labeled | BDE | EO-5403 (ng/mL) | EO-5403-10X-1.2 (ng/mL) |
|---|-----|--------------------|----------------------------|
| 4-MonoBDE ($^{13}\text{C}_{12}$, 99%) | 3 | 100 | 1000 |
| 4,4'-DiBDE ($^{13}\text{C}_{12}$, 99%) | 15 | 100 | 1000 |
| 2,4,4'-TriBDE ($^{13}\text{C}_{12}$, 99%) | 28 | 100 | 1000 |
| 2,2',4,4'-TetraBDE ($^{13}\text{C}_{12}$, 99%) | 47 | 100 | 1000 |
| 2,2',4,4',5-PentaBDE ($^{13}\text{C}_{12}$, 99%) | 99 | 100 | 1000 |
| 2,2',4,4',5,5'-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 153 | 200 | 2000 |
| 2,2',4,4',5,6'-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 154 | 200 | 2000 |
| 2,2',3,4,4',5',6-HeptaBDE ($^{13}\text{C}_{12}$, 99%) | 183 | 200 | 2000 |
| 2,2',3,4,4',5,6,6'-OctaBDE ($^{13}\text{C}_{12}$, 99%) | 204 | 200 | 2000 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE ($^{13}\text{C}_{12}$, 99%) | 207 | 500 | 5000 |
| DecaBDE ($^{13}\text{C}_{12}$, 99%) | 209 | 500 | 5000 |

| | | |
|---------|-------------------------|------------------|
| EO-5404 | RoHS PBDE Syringe Spike | 1.2 mL in nonane |
|---------|-------------------------|------------------|

| Labeled | BDE | (ng/mL) |
|--|-----|---------|
| 2,2',3,4,4',5'-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 138 | 200 |

| | | |
|---------|----------------------------|------------------|
| EO-5405 | RoHS PBDE Native PAR Spike | 1.2 mL in nonane |
|---------|----------------------------|------------------|

| Unlabeled | BDE | (ng/mL) |
|-------------------------------|-----|---------|
| 4-MonoBDE | 3 | 1000 |
| 2,4-DiBDE | 7 | 1000 |
| 4,4'-DiBDE | 15 | 1000 |
| 2,2',4-TriBDE | 17 | 1000 |
| 2,4,4'-TriBDE | 28 | 1000 |
| 2,2',4,4'-TetraBDE | 47 | 1000 |
| 2,2',4,5'-TetraBDE | 49 | 1000 |
| 2,3',4,4'-TetraBDE | 66 | 1000 |
| 2,3',4',6-TetraBDE | 71 | 1000 |
| 3,3',4,4'-TetraBDE | 77 | 1000 |
| 2,2',3,4,4'-PentaBDE | 85 | 1000 |
| 2,2',4,4',5-PentaBDE | 99 | 1000 |
| 2,2',4,4',6-PentaBDE | 100 | 1000 |
| 2,3',4,4',6-PentaBDE | 119 | 1000 |
| 3,3',4,4',5-PentaBDE | 126 | 1000 |
| 2,2',3,4,4',5'-HexaBDE | 138 | 2000 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 2000 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 2000 |
| 2,2',4,4',6,6'-HexaBDE | 155 | 2000 |
| 2,3,4,4',5,6-HexaBDE | 166 | 2000 |
| 2,2',3,4,4',5,6-HeptaBDE | 181 | 2000 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 2000 |
| 2,3,3',4,4',5,6-HeptaBDE | 190 | 2000 |
| 2,2',3,4,4',5,5',6-OctaBDE | 203 | 2000 |
| 2,3,3',4,4',5,5',6-OctaBDE | 205 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaBDE | 206 | 5000 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE | 207 | 5000 |
| DecaBDE | 209 | 5000 |

RoHS Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EO-5425 | RoHS Screening PBDE Calibration Solutions [CS1-CS3] | Set of 3 x 0.2 mL in nonane |
| EO-5425-CS1 | RoHS Screening PBDE Calibration Solution [CS1] | 0.2 mL in nonane |
| EO-5425-CS2 | RoHS Screening PBDE Calibration Solution [CS2] | 0.2 mL in nonane |
| EO-5425-CS3 | RoHS Screening PBDE Calibration Solution [CS3] | 0.2 mL in nonane |

| <i>All concentrations are in ng/mL (ppb)</i> | | | | | |
|--|-----|------|------|--------|--|
| Unlabeled | BDE | CS1 | CS2 | CS3 | |
| 4,4'-DiBDE | 15 | 20 | 100 | 1000 | |
| 2,4,4'-TriBDE | 28 | 20 | 100 | 1000 | |
| 2,2',4,4'-TetraBDE | 47 | 20 | 100 | 1000 | |
| 2,2',4,4',5-PentaBDE | 99 | 20 | 100 | 1000 | |
| 2,2',4,4',5,5'-HexaBDE | 153 | 20 | 100 | 1000 | |
| 2,2',4,4',5,6'-HexaBDE | 154 | 20 | 100 | 1000 | |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 20 | 100 | 1000 | |
| 2,2',3,3',4,4',6,6'-OctaBDE | 197 | 50 | 250 | 2500 | |
| 2,2',3,3',4,4',5,5',6-NonaBDE | 206 | 100 | 500 | 5000 | |
| 2,2',3,3',4,4',5,6,6'-NonaBDE | 207 | 50 | 250 | 2500 | |
| 2,2',3,3',4,5,5',6,6'-NonaBDE | 208 | 50 | 250 | 2500 | |
| DecaBDE | 209 | 500 | 2500 | 25,000 | |
| Cleanup | | | | | |
| 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 100 | 100 | 100 | |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 | 100 | 100 | |
| 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 100 | 100 | 100 | |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 100 | 100 | 100 | |
| 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 100 | 100 | 100 | |
| 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 100 | 100 | 100 | |
| 2,2',3,3',4,4',6,6'-OctaBDE (¹³ C ₁₂ , 99%) | 197 | 250 | 250 | 250 | |
| 2,2',3,3',4,4',5,5',6-NonaBDE (¹³ C ₁₂ , 99%) | 206 | 500 | 500 | 500 | |
| 2,2',3,3',4,5,5',6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 208 | 250 | 250 | 250 | |
| DecaBDE (¹³ C ₁₂ , 99%) | 209 | 2500 | 2500 | 2500 | |
| Syringe | | | | | |
| 2,2',3,3',4,4',5,6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 207 | 250 | 250 | 250 | |

RoHS Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | | Amount |
|-------------|--|-----|------------------|
| EO-5426 | RoHS Screening PBDE Cleanup Spike | | 1.2 mL in nonane |
| | Labeled | BDE | (ng/mL) |
| | 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 1000 |
| | 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 1000 |
| | 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 1000 |
| | 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 1000 |
| | 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 1000 |
| | 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 1000 |
| | 2,2',3,4,4',5'-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 1000 |
| | 2,2',3,3',4,4',6,6'-OctaBDE (¹³ C ₁₂ , 99%) | 197 | 2500 |
| | 2,2',3,3',4,4',5,5'-NonaBDE (¹³ C ₁₂ , 99%) | 206 | 5000 |
| | 2,2',3,3',4,5,5',6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 208 | 2500 |
| | DecaBDE (¹³ C ₁₂ , 99%) | 209 | 25,000 |
| EO-5427 | RoHS Screening PBDE Syringe Spike | | 1.2 mL in nonane |
| | Labeled | BDE | (ng/mL) |
| | 2,2',3,3',4,4',5,6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 207 | 2500 |
| EO-5428 | RoHS Screening PBDE Native PAR Spike | | 1.2 mL in nonane |
| | Unlabeled | BDE | (ng/mL) |
| | 4,4'-DiBDE | 15 | 1000 |
| | 2,4,4'-TriBDE | 28 | 1000 |
| | 2,2',4,4'-TetraBDE | 47 | 1000 |
| | 2,2',4,4',5-PentaBDE | 99 | 1000 |
| | 2,2',4,4',5,5'-HexaBDE | 153 | 1000 |
| | 2,2',4,4',5,6'-HexaBDE | 154 | 1000 |
| | 2,2',3,4,4',5'-HeptaBDE | 183 | 1000 |
| | 2,2',3,3',4,4',6,6'-OctaBDE | 197 | 2500 |
| | 2,2',3,3',4,4',5,5'-NonaBDE | 206 | 5000 |
| | 2,2',3,3',4,4',5,6,6'-NonaBDE | 207 | 2500 |
| | 2,2',3,3',4,5,5',6,6'-NonaBDE | 208 | 2500 |
| | DecaBDE | 209 | 25,000 |

Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|-----------------------------|
| EO-5104 | Brominated Diphenyl Ether Calibration Solutions [CS1-CS6] | Set of 6 x 0.2 mL in nonane |
| EO-5104-CS1 | Brominated Diphenyl Ether Calibration Solution [CS1] | 0.2 mL in nonane |
| EO-5104-CS2 | Brominated Diphenyl Ether Calibration Solution [CS2] | 0.2 mL in nonane |
| EO-5104-CS3 | Brominated Diphenyl Ether Calibration Solution [CS3] | 0.2 mL in nonane |
| EO-5104-CS4 | Brominated Diphenyl Ether Calibration Solution [CS4] | 0.2 mL in nonane |
| EO-5104-CS5 | Brominated Diphenyl Ether Calibration Solution [CS5] | 0.2 mL in nonane |
| EO-5104-CS6 | Brominated Diphenyl Ether Calibration Solution [CS6] | 0.2 mL in nonane |

| All concentrations are in ng/mL (ppb) | | | | | | | |
|---|-----|-----|-----|------|------|-----|------|
| Unlabeled | BDE | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
| 2-MonoBDE | 1 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3-MonoBDE | 2 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 4-MonoBDE | 3 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4-DiBDE | 7 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4'-DiBDE | 8 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,6-DiBDE | 10 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,3'-DiBDE | 11 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,4-DiBDE | 12 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,4'-DiBDE | 13 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 4,4'-DiBDE | 15 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,2',4-TriBDE | 17 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,3',4-TriBDE | 25 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4,4'-TriBDE | 28 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4,6-TriBDE | 30 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4',6-TriBDE | 32 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2',3,4-TriBDE | 33 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,3',4-TriBDE | 35 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,4,4'-TriBDE | 37 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,2',4,4'-TetraBDE | 47 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,2',4,5'-TetraBDE | 49 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,3',4,4'-TetraBDE | 66 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,3',4',6-TetraBDE | 71 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,4,4',6-TetraBDE | 75 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 3,3',4,4'-TetraBDE | 77 | 0.2 | 1 | 5 | 25 | 100 | 500 |
| 2,2',3,4,4'-PentabDE | 85 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,2',4,4',5-PentabDE | 99 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,2',4,4',6-PentabDE | 100 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,3,4,5,6-PentaBDE | 116 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,3',4,4',5-PentabDE | 118 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,3',4,4',6-PentabDE | 119 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 3,3',4,4',5-PentabDE | 126 | 0.3 | 1.5 | 7.5 | 37.5 | 150 | 750 |
| 2,2',3,4,4',5'-HexaBDE | 138 | 0.6 | 3 | 15 | 75 | 300 | 1500 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 0.4 | 2 | 10 | 50 | 200 | 1000 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 0.4 | 2 | 10 | 50 | 200 | 1000 |
| 2,2',4,4',6,6'-HexaBDE | 155 | 0.4 | 2 | 10 | 50 | 200 | 1000 |
| 2,3,4,4',5,6-HexaBDE | 166 | 0.4 | 2 | 10 | 50 | 200 | 1000 |
| 2,2',3,4,4',5,6-HeptaBDE | 181 | 0.5 | 2.5 | 12.5 | 62.5 | 250 | 1250 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 0.5 | 2.5 | 12.5 | 62.5 | 250 | 1250 |
| 2,3,3',4,4',5,6-HeptaBDE | 190 | 0.5 | 2.5 | 12.5 | 62.5 | 250 | 1250 |
| Surrogate | | | | | | | |
| 4-MonoBDE (¹³ C ₁₂ , 99%) | 3 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5-PentabDE (¹³ C ₁₂ , 99%) | 99 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',6-PentabDE (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5-PentabDE (¹³ C ₁₂ , 99%) | 118 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 200 | 200 | 200 | 200 | 200 | 200 |
| 2,2',3,4,4',5,6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 250 | 250 | 250 | 250 | 250 | 250 |
| Performance | | | | | | | |
| 3,3',4,4',4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4',5-PentabDE (¹³ C ₁₂ , 99%) | 126 | 150 | 150 | 150 | 150 | 150 | 150 |

Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|-----------------|--|------------------|
| EO-5100 | Polybrominated Diphenyl Ether Surrogate Spiking Solution | 1.2 mL in nonane |
| EO-5100-10X-0.5 | Polybrominated Diphenyl Ether Surrogate Spiking Solution (10X stock) | 0.5 mL in nonane |

| Labeled | BDE | EO-5100 (ng/mL) | EO-5100-10X-0.5 (ng/mL) |
|--|-----|--------------------|----------------------------|
| 4-MonoBDE (¹³ C ₁₂ , 99%) | 3 | 100 | 1000 |
| 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 100 | 1000 |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 | 1000 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 | 1000 |
| 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 150 | 1500 |
| 2,2',4,4',6-PentaBDE (¹³ C ₁₂ , 99%) | 100 | 150 | 1500 |
| 2,3',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 118 | 150 | 1500 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 200 | 2000 |
| 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 250 | 2500 |

| | | |
|-----------------|---|------------------|
| EO-5101 | Polybrominated Diphenyl Ether Performance Standard Solution | 1.2 mL in nonane |
| EO-5101-10X-1.2 | Polybrominated Diphenyl Ether Performance Standard Solution (10X stock) | 1.2 mL in nonane |

| Labeled | BDE | EO-5101 (ng/mL) | EO-5101-10X-1.2 (ng/mL) |
|---|-----|--------------------|----------------------------|
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 100 | 1000 |
| 3,3',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 126 | 150 | 1500 |

| | | |
|--------------------|---|------------------|
| NEW EO-5507 | Polybrominated Diphenyl Ether Injection Internal Standard | 1.2 mL in nonane |
|--------------------|---|------------------|

| Labeled | BDE | (ng/mL) |
|---|-----|---------|
| 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 200 |
| 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 200 |

| | | |
|--------------------|--|------------------|
| NEW EO-5509 | Polybrominated Diphenyl Ether Native Injection Internal Standard | 1.2 mL in nonane |
|--------------------|--|------------------|

| Unlabeled | BDE | (ng/mL) |
|------------------------|-----|---------|
| 2,2',4,4',5-PentaBDE | 99 | 200 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 200 |

| | | |
|--------------------|--|------------------|
| NEW EO-5506 | Polybrominated Diphenyl Ether Surrogate Standard | 1.2 mL in nonane |
|--------------------|--|------------------|

| Labeled | BDE | (ng/mL) |
|--|-----|---------|
| 2,2',4,4',6-PentaBDE (¹³ C ₁₂ , 99%) | 100 | 200 |
| 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 200 |

| | | |
|--------------------|---|------------------|
| NEW EO-5508 | Polybrominated Diphenyl Ether Native Surrogate Standard | 1.2 mL in nonane |
|--------------------|---|------------------|

| Unlabeled | BDE | (ng/mL) |
|---------------------------|-----|---------|
| 2,2',4,4',6-PentaBDE | 100 | 200 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 200 |

Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| EO-5099 | Polybrominated Diphenyl Ether Analytical Standard Solution | 1.2 mL in nonane |
| Unlabeled | | |
| 2-MonoBDE | 1 | 100 |
| 3-MonoBDE | 2 | 100 |
| 4-MonoBDE | 3 | 100 |
| 2,4-DiBDE | 7 | 100 |
| 2,4'-DiBDE | 8 | 100 |
| 2,6-DiBDE | 10 | 100 |
| 3,3'-DiBDE | 11 | 100 |
| 3,4-DiBDE | 12 | 100 |
| 3,4'-DiBDE | 13 | 100 |
| 4,4'-DiBDE | 15 | 100 |
| 2,2',4-TriBDE | 17 | 100 |
| 2,3',4-TriBDE | 25 | 100 |
| 2,4,4'-TriBDE | 28 | 100 |
| 2,4,6-TriBDE | 30 | 100 |
| 2,4',6-TriBDE | 32 | 100 |
| 2',3,4-TriBDE | 33 | 100 |
| 3,3',4-TriBDE | 35 | 100 |
| 3,4,4'-TriBDE | 37 | 100 |
| 2,2',4,4'-TetraBDE | 47 | 100 |
| 2,2',4,5'-TetraBDE | 49 | 100 |
| 2,3',4,4'-TetraBDE | 66 | 100 |
| 2,3',4',6-TetraBDE | 71 | 100 |
| 2,4,4',6-TetraBDE | 75 | 100 |
| 3,3',4,4'-TetraBDE | 77 | 100 |
| 2,2',3,4,4'-PentaBDE | 85 | 150 |
| 2,2',4,4',5-PentaBDE | 99 | 150 |
| 2,2',4,4',6-PentaBDE | 100 | 150 |
| 2,3,4,5,6-PentaBDE | 116 | 150 |
| 2,3',4,4',5-PentaBDE | 118 | 150 |
| 2,3',4,4',6-PentaBDE | 119 | 150 |
| 3,3',4,4',5-PentaBDE | 126 | 150 |
| 2,2',3,4,4',5-HexaBDE | 138 | 300 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 200 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 200 |
| 2,2',4,4',6,6'-HexaBDE | 155 | 200 |
| 2,3,4,4',5,6-HexaBDE | 166 | 200 |
| 2,2',3,4,4',5,6-HeptaBDE | 181 | 250 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 250 |
| 2,3,3',4,4',5,6-HeptaBDE | 190 | 250 |
| Labeled | | |
| 4-MonoBDE (¹³ C ₁₂ , 99%) | 3 | 100 |
| 4,4'-DiBDE (¹³ C ₁₂ , 99%) | 15 | 100 |
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 100 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 100 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 100 |
| 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 150 |
| 2,2',4,4',6-PentaBDE (¹³ C ₁₂ , 99%) | 100 | 150 |
| 2,3',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 118 | 150 |
| 3,3',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 126 | 150 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 200 |
| 2,2',3,4,4',5,6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 250 |

Brominated Diphenyl Ether (BDE) Standard Mixtures

| Catalog No. | Compound | | | Amount |
|-------------|--|-----|---------|---------------------------|
| EO-5103 | Polybrominated Diphenyl Ether Predominant Congener Mixture | | | 1.2 mL in nonane |
| | Unlabeled | BDE | (ng/mL) | Unlabeled |
| | 2,2',4-TriBDE | 17 | 2500 | 2,2',4,4',6-PentaBDE |
| | 2,4,4'-TriBDE | 28 | 2500 | 2,2',3,4,4',5'-HexaBDE |
| | 2,2',4,4'-TetraBDE | 47 | 2500 | 2,2',4,4',5,5'-HexaBDE |
| | 2,3',4,4'-TetraBDE | 66 | 2500 | 2,2',4,4',5,6'-HexaBDE |
| | 2,3',4',6-TetraBDE | 71 | 2500 | 2,2',3,4,4',5',6-HeptaBDE |
| | 2,2',3,4,4'-PentaBDE | 85 | 2500 | 2,3,3',4,4',5,6-HeptaBDE |
| | 2,2',4,4',5-PentaBDE | 99 | 2500 | DecaBDE |

| | | |
|------------------|---|------------------|
| EO-5113 | Polybrominated Diphenyl Ether PAR Solution | 0.5 mL in nonane |
| EO-5113-7.5X-0.5 | Polybrominated Diphenyl Ether PAR Solution (7.5X stock) | 0.5 mL in nonane |

| | | EO-5113 (ng/mL) | EO-5113-7.5X-0.5 (ng/mL) |
|--|---------------------------|--------------------|-----------------------------|
| | Unlabeled | BDE | |
| | 2-MonoBDE | 1 | 100 |
| | 3-MonoBDE | 2 | 100 |
| | 4-MonoBDE | 3 | 100 |
| | 2,4-DiBDE | 7 | 100 |
| | 2,4'-DiBDE | 8 | 100 |
| | 2,6-DiBDE | 10 | 100 |
| | 3,3'-DiBDE | 11 | 100 |
| | 3,4-DiBDE | 12 | 100 |
| | 3,4'-DiBDE | 13 | 100 |
| | 4,4'-DiBDE | 15 | 100 |
| | 2,2',4-TriBDE | 17 | 100 |
| | 2,3',4-TriBDE | 25 | 100 |
| | 2,4,4'-TriBDE | 28 | 100 |
| | 2,4,6-TriBDE | 30 | 100 |
| | 2,4',6-TriBDE | 32 | 100 |
| | 2',3,4-TriBDE | 33 | 100 |
| | 3,3',4-TriBDE | 35 | 100 |
| | 3,4,4'-TriBDE | 37 | 100 |
| | 2,2',4,4'-TetraBDE | 47 | 100 |
| | 2,2',4,5'-TetraBDE | 49 | 100 |
| | 2,3',4,4'-TetraBDE | 66 | 100 |
| | 2,3',4',6-TetraBDE | 71 | 100 |
| | 2,4,4',6-TetraBDE | 75 | 100 |
| | 3,3',4,4'-TetraBDE | 77 | 100 |
| | 2,2',3,4,4'-PentaBDE | 85 | 150 |
| | 2,2',4,4',5-PentaBDE | 99 | 150 |
| | 2,2',4,4',6-PentaBDE | 100 | 150 |
| | 2,3,4,5,6-PentaBDE | 116 | 150 |
| | 2,3',4,4',5-PentaBDE | 118 | 150 |
| | 2,3',4,4',6-PentaBDE | 119 | 150 |
| | 3,3',4,4',5-PentaBDE | 126 | 150 |
| | 2,2',3,4,4',5'-HexaBDE | 138 | 300 |
| | 2,2',4,4',5,5'-HexaBDE | 153 | 200 |
| | 2,2',4,4',5,6'-HexaBDE | 154 | 200 |
| | 2,2',4,4',6,6'-HexaBDE | 155 | 200 |
| | 2,3,4,4',5,6-HexaBDE | 166 | 200 |
| | 2,2',3,4,4',5,6-HeptaBDE | 181 | 250 |
| | 2,2',3,4,4',5',6-HeptaBDE | 183 | 250 |
| | 2,3,3',4,4',5,6-HeptaBDE | 190 | 250 |

Brominated Diphenyl Ether (BDE) Technical Mixtures

| Catalog No. | Compound | Concentration | Amount |
|-------------|--|----------------------|--------|
| EO-4958-1.2 | Pentabromodiphenyl Ether Technical Mix (Bromkal™ 70-5) | 50 µg/mL in methanol | 1.2 mL |
| EO-5031 | Pentabromodiphenyl Ether Technical Mix (DE-71™) | 50 µg/mL in methanol | 1.2 mL |
| EO-5030 | Octabromodiphenyl Ether Technical Mix (DE-79™) | 50 µg/mL in methanol | 1.2 mL |
| EO-5060 | Decabromodiphenyl Ether Technical Mix (Saytex® 102E) | 10 µg/mL in methanol | 10 mL |

Bromkal is a trademark of Chemische Fabrik Kalk GmbH. DE-71 is a trademark of Chemtura (Great Lakes Chemical Company). DE-79 is a trademark of Chemtura (Great Lakes Chemical Company). Saytex is a registered trademark of Albemarle Corporation.

Brominated Diphenyl Ether (BDE) Metabolite Standards

| Catalog No. | Compound | Concentration | Amount |
|----------------------------|---|---------------------|--------|
| OH-BDE | | | |
| OHBDE-5190-1.2 | 6-Hydroxy-2,2',4,4'-tetrabromodiphenyl ether (ring- ¹³ C ₁₂ , 99%) CP 92% | 50 µg/mL in toluene | 1.2 mL |
| OHBDE-5206-1.2 | 6-Hydroxy-2,2',4,4'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| OHBDE-5191-1.2 | 2-Hydroxy-2',4,4',5,6-pentabromodiphenyl ether (ring- ¹³ C ₁₂ , 99%) CP 94% | 50 µg/mL in toluene | 1.2 mL |
| OHBDE-5212-1.2 | 4'-Hydroxy-2,2',4,5'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| OHBDE-5228-1.2 | 6-Hydroxy-2,2',4,4',5-pentabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| MeO-BDE | | | |
| MEOBDE-5153-1.2 | 2'-Methoxy-2,3',4,5'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5207-1.2 | 3-Methoxy-2,2',4,4'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5203-1.2 | 4-Methoxy-2,2',3,4'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5211-1.2 | 4'-Methoxy-2,2',4,5'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5209-1.2 | 5-Methoxy-2,2',4,4'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5260-1.2 | 6-Methoxy-2,2',4,4'-tetrabromodiphenyl ether (ring- ¹³ C ₁₂ , 99%) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5205-1.2 | 6-Methoxy-2,2',4,4'-tetrabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| <i>NEW</i> MEOBDE-5227-1.2 | 6-Methoxy-2,2',4,4',5-pentabromodiphenyl ether (unlabeled) | 50 µg/mL in nonane | 1.2 mL |

Tetrabromobisphenol A (TBBPA) and Tetrachlorobisphenol A Standards

| Catalog No. | Compound | Concentration | Amount |
|---------------------------|--|----------------------|--------|
| <i>NEW</i> CLM-9374-1.2 | Dimethyl tetrabromobisphenol A (ring- ¹³ C ₁₂ , 99%) | 100 µg/mL in toluene | 1.2 mL |
| ULM-6236-1.2 | Dimethyl tetrabromobisphenol A (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| CLM-4694-1.2 | Tetrabromobisphenol A (TBBPA) (ring- ¹³ C ₁₂ , 99%) | 50 µg/mL in methanol | 1.2 mL |
| <i>NEW</i> CLM-4694-T-1.2 | Tetrabromobisphenol A (TBBPA) (ring- ¹³ C ₁₂ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| ULM-8734-1.2 | Tetrabromobisphenol A (TBBPA) (unlabeled) | 50 µg/mL in methanol | 1.2 mL |
| <i>NEW</i> ULM-8734-T-1.2 | Tetrabromobisphenol A (TBBPA) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| CLM-8006-1.2 | Tetrachlorobisphenol A (ring- ¹³ C ₁₂ , 99%) | 50 µg/mL in methanol | 1.2 mL |
| ULM-7606-1.2 | Tetrachlorobisphenol A (unlabeled) | 50 µg/mL in methanol | 1.2 mL |

Hexabromocyclododecane (HBCD) Standards

| Catalog No. | Compound | Concentration | Amount |
|--------------|---|---------------------|--------|
| CLM-7922-0.5 | α-Hexabromocyclododecane (HBCD) (¹³ C ₁₂ , 99%) | 50 µg/mL in toluene | 0.5 mL |
| ULM-4834-1.2 | α-Hexabromocyclododecane (HBCD) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| CLM-7923-1.2 | β-Hexabromocyclododecane (HBCD) (¹³ C ₁₂ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| ULM-4835-1.2 | β-Hexabromocyclododecane (HBCD) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| CLM-7924-1.2 | γ-Hexabromocyclododecane (HBCD) (¹³ C ₁₂ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| ULM-4836-1.2 | γ-Hexabromocyclododecane (HBCD) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| CLM-7102-1.2 | Hexabromocyclododecane (HBCD) (¹³ C ₁₂ , 99%) unequal mixture of α, β, and γ isomers | 50 µg/mL in toluene | 1.2 mL |

Dechlorane Standards

| Catalog No. | Compound | Concentration | Amount |
|--------------------|---|----------------------|--------|
| CLM-8569-1.2 | Dechlorane Plus <i>syn</i> (bis-cyclopentene- ¹³ C ₁₀ , 99%) | 100 µg/mL in nonane | 1.2 mL |
| CLM-8569-T-1.2 | | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-7886-1.2 | Dechlorane Plus <i>syn</i> (unlabeled) | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-7886-T-1.2 | | 100 µg/mL in toluene | 1.2 mL |
| CLM-8588-1.2 | Dechlorane Plus <i>anti</i> (bis-cyclopentene- ¹³ C ₁₀ , 99%) | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-8588-T-1.2 | | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-7887-1.2 | Dechlorane Plus <i>anti</i> (unlabeled) | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-7887-T-1.2 | | 100 µg/mL in toluene | 1.2 mL |
| ULM-7777-1.2 | Dechlorane Plus Technical Product (unlabeled) mix of <i>syn/anti</i> isomers | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9282-1.2 | Dechlorane 602 (¹³ C ₁₀ , 99%) | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9283-1.2 | Dechlorane 602 (unlabeled) | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9621-1.2 | Dechlorane 603 (unlabeled) | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9622-1.2 | Dechlorane 604 (Component A) (unlabeled) CP >95% | 100 µg/mL in nonane | 1.2 mL |

Polybrominated Biphenyl (PBB) Standards

| Catalog No. | Compound | PBB | Concentration | Amount |
|-------------|---|-----|-----------------------|--------|
| EB-5055 | 3,3',4,4'-Tetrabromobiphenyl (¹³ C ₁₂ , 99%) | 77 | 40 µg/mL in nonane | 3 mL |
| PBB-77-CS | 3,3',4,4'-Tetrabromobiphenyl Certified Standard (unlabeled) | 77 | 100 µg/mL in isoctane | 1.2 mL |
| EB-5056 | 3,3',4,4',5-Pentabromobiphenyl (¹³ C ₁₂ , 99%) | 126 | 40 µg/mL in nonane | 3 mL |
| PBB-126 | 3,3',4,4',5-Pentabromobiphenyl (unlabeled) | 126 | 100 µg/mL in isoctane | 1.2 mL |
| EB-5162 | 2,2',4,4',5,5'-Hexabromobiphenyl (¹³ C ₁₂ , 99%) | 153 | 40 µg/mL in nonane | 3 mL |
| PBB-153-CS | 2,2',4,4',5,5'-Hexabromobiphenyl Certified Standard (unlabeled) | 153 | 100 µg/mL in isoctane | 1.2 mL |
| EB-5106 | 2,3,3',4,4',5'-Hexabromobiphenyl (¹³ C ₁₂ , 99%) | 157 | 40 µg/mL in nonane | 3 mL |
| PBB-157-CS | 2,3,3',4,4',5'-Hexabromobiphenyl Certified Standard (unlabeled) | 157 | 100 µg/mL in isoctane | 1.2 mL |
| EB-5439 | Decabromobiphenyl (¹³ C ₁₂ , 99%) | 209 | 40 µg/mL in nonane | 3 mL |
| PBB-209-CS | Decabromobiphenyl Certified Standard (unlabeled) | 209 | 100 µg/mL in isoctane | 1.2 mL |

Other Flame-Retardant Standards

| Catalog No. | Compound | Concentration | Amount |
|------------------|--|----------------------|------------|
| ULM-7375-2X1.2 | 1,2-Bis(pentabromophenyl) ethane (DBDPE) (unlabeled) CP 96% | 25 µg/mL in toluene | 2 × 1.2 mL |
| ULM-7595-1.2 | 1,2-Bis(2,4,6-tribromophenoxy) ethane (BTBPE) (unlabeled) | 50 µg/mL in nonane | 1.2 mL |
| NEW CLM-1921-1.2 | Hexabromobenzene (¹³ C ₆ , 99%) | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-7607-1.2 | Hexabromobenzene (unlabeled) | 100 µg/mL in toluene | 1.2 mL |
| NEW CLM-9535-1.2 | Tris(2,4,6-tribromophenoxy)-1,3,5-triazine (TTBP-TAZ) (¹³ C ₁₈ , 99%) | 50 µg/mL in dioxane | 1.2 mL |
| NEW ULM-9506-1.2 | Tris(2,4,6-tribromophenoxy)-1,3,5-triazine (TTBP-TAZ) (unlabeled) | 50 µg/mL in dioxane | 1.2 mL |
| NEW DLM-9945-1.2 | TBB (2-ethylhexyl-2,3,4,5-tetrabromobenzoate) (D ₁₇ , 98%) | 50 µg/mL in toluene | Inquire |
| NEW ULM-9944-1.2 | TBB (2-ethylhexyl-2,3,4,5-tetrabromobenzoate) (unlabeled) | 50 µg/mL in toluene | Inquire |
| NEW DLM-9947-1.2 | TBPH (bis(2-ethylhexyl)tetrabromophthalate) (D ₃₄ , 98%) | 50 µg/mL in toluene | Inquire |
| NEW ULM-9946-1.2 | TBPH (bis(2-ethylhexyl)tetrabromophthalate) (unlabeled) | 50 µg/mL in toluene | Inquire |

Brominated Flame-Retardant (BFR) Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|--|------------------------------|
| EO-5319-A | CDC BFR Calibration Standards [CS1-CS10] | Set of 10 x 0.5 mL in nonane |
| EO-5319-A-CS1 | CDC BFR Calibration Standard [CS1] | 0.5 mL in nonane |
| EO-5319-A-CS2 | CDC BFR Calibration Standard [CS2] | 0.5 mL in nonane |
| EO-5319-A-CS3 | CDC BFR Calibration Standard [CS3] | 0.5 mL in nonane |
| EO-5319-A-CS4 | CDC BFR Calibration Standard [CS4] | 0.5 mL in nonane |
| EO-5319-A-CS5 | CDC BFR Calibration Standard [CS5] | 0.5 mL in nonane |
| EO-5319-A-CS6 | CDC BFR Calibration Standard [CS6] | 0.5 mL in nonane |
| EO-5319-A-CS7 | CDC BFR Calibration Standard [CS7] | 0.5 mL in nonane |
| EO-5319-A-CS8 | CDC BFR Calibration Standard [CS8] | 0.5 mL in nonane |
| EO-5319-A-CS9 | CDC BFR Calibration Standard [CS9] | 0.5 mL in nonane |
| EO-5319-A-CS10 | CDC BFR Calibration Standard [CS10] | 0.5 mL in nonane |

All concentrations are in ng/mL (ppb)

| Unlabeled | Congener | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 | CS10 |
|---------------------------------------|----------|-----|-----|-----|-----|-----|-----|-----|-----|------|------|
| 2,2',4-TriBDE | 17 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,4,4'-TriBDE | 28 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4'-TetraBDE | 47 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,3',4,4'-TetraBDE | 66 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4',6-PentaBDE | 100 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4',5-PentaBDE | 99 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,4,4'-PentaBDE | 85 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4',5,6'-HexaBDE | 154 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4',5,5'-HexaBDE | 153 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,4,4',5',6-HeptaBDE | 183 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,3',4,4',5,6'-OctaBDE | 196 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,3',4,4',6,6'-OctaBDE | 197 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,4,4',5,5',6-OctaBDE | 203 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,3',4,4',5,5',6-NonaBDE | 206 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE | 207 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',3,3',4,5,5',6,6'-NonaBDE | 208 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| DecaBDE | 209 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 2,2',4,4',5,5'-HexaBB | PBB-153 | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| Hexabromobenzene | | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 1,2-Bis(pentabromophenyl) ethane | | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| 1,2-Bis(2,4,6-tribromophenoxy) ethane | | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |
| γ -Hexabromobenzene | | 0.2 | 0.5 | 1 | 5 | 10 | 50 | 100 | 500 | 1000 | 2000 |

| Labeled | | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
|---|---------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| 1,2,3,4-TetraCDD ($^{13}\text{C}_6$, 99%) | | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6,6'-NonaCB ($^{13}\text{C}_{12}$, 99%) | 208 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-TriBDE ($^{13}\text{C}_{12}$, 99%) | 28 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4'-TetraBDE ($^{13}\text{C}_{12}$, 99%) | 47 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 3,3',4,4'-TetraBDE ($^{13}\text{C}_{12}$, 99%) | 77 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4',6-PentaBDE ($^{13}\text{C}_{12}$, 99%) | 100 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4',5-PentaBDE ($^{13}\text{C}_{12}$, 99%) | 99 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4',5,6'-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 154 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',4,4',5,5'-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 153 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-HexaBDE ($^{13}\text{C}_{12}$, 99%) | 139 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',5'-HeptaBDE ($^{13}\text{C}_{12}$, 99%) | 183 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',6,6'-OctaBDE ($^{13}\text{C}_{12}$, 99%) | 197 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',5,5',6-OctaBDE ($^{13}\text{C}_{12}$, 99%) | 203 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5,5',6-NonaBDE ($^{13}\text{C}_{12}$, 99%) | 206 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE ($^{13}\text{C}_{12}$, 99%) | 207 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,3',4,4',5,5',6-NonaBDE ($^{13}\text{C}_{12}$, 99%) | 208 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| DecaBDE ($^{13}\text{C}_{12}$, 99%) | 209 | 500 | 500 | 500 | 500 | 500 | 500 | 500 | 500 | 500 | 500 |
| 2,2',4,4',5,5'-HexaBB ($^{13}\text{C}_{12}$, 99%) | PBB-153 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Hexabromobenzene ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 1,2-Bis(pentabromophenyl) ethane ($^{13}\text{C}_{14}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 1,2-Bis(2,4,6-tribromophenoxy) ethane ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| γ -Hexabromocyclododecane ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |

Brominated Flame-Retardant (BFR) Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------|--------------------------|--------------------------|
| EO-5320-A | CDC BFR Spiking Standard | 10 mL in methanol |
| EO-5320-200X-1.2 | CDC BFR Spiking Standard | 1.2 mL in nonane/toluene |

| Labeled | Congener | EC-5320-A (ng/mL) | EO-5320-200X-1.2 (ng/mL) |
|--|----------|----------------------|-----------------------------|
| 2,4,4'-TriBDE (¹³ C ₁₂ , 99%) | 28 | 7.5 | 1500 |
| 2,2',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 47 | 7.5 | 1500 |
| 2,2',4,4',6-PentaBDE (¹³ C ₁₂ , 99%) | 100 | 7.5 | 1500 |
| 2,2',4,4',5-PentaBDE (¹³ C ₁₂ , 99%) | 99 | 7.5 | 1500 |
| 2,2',4,4',5,6'-HexaBDE (¹³ C ₁₂ , 99%) | 154 | 7.5 | 1500 |
| 2,2',4,4',5,5'-HexaBDE (¹³ C ₁₂ , 99%) | 153 | 7.5 | 1500 |
| 2,2',3,4,4',5',6-HeptaBDE (¹³ C ₁₂ , 99%) | 183 | 7.5 | 1500 |
| 2,2',3,3',4,4',6,6'-OctaBDE (¹³ C ₁₂ , 99%) | 197 | 7.5 | 1500 |
| 2,2',3,4,4',5,5',6-OctaBDE (¹³ C ₁₂ , 99%) | 203 | 7.5 | 1500 |
| 2,2',3,3',4,4',5,5',6-NonaBDE (¹³ C ₁₂ , 99%) | 206 | 7.5 | 1500 |
| 2,2',3,3',4,4',5,6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 207 | 7.5 | 1500 |
| 2,2',3,3',4,5,5',6,6'-NonaBDE (¹³ C ₁₂ , 99%) | 208 | 7.5 | 1500 |
| DecaBDE (¹³ C ₁₂ , 99%) | 209 | 50 | 10,000 |
| 2,2',4,4',5,5'-HexaBB (¹³ C ₁₂ , 99%) | PBB-153 | 7.5 | 1500 |
| Hexabromobenzene (¹³ C ₆ , 99%) | | 7.5 | 1500 |
| 1,2-Bis(pentabromophenyl) ethane (¹³ C ₁₄ , 99%) | | 7.5 | 1500 |
| 1,2-Bis(2,4,6-tribromophenoxy) ethane (¹³ C ₁₂ , 99%) | | 7.5 | 1500 |
| γ-Hexabromocyclododecane (¹³ C ₁₂ , 99%) | | 7.5 | 1500 |

| | | |
|---------|-------------------------------|--|
| EO-5169 | BFR Recovery Spiking Solution | 10 mL in 88% hexane/ 2% dodecane/10% nonane |
|---------|-------------------------------|--|

| Labeled | Congener | (ng/mL) |
|---|----------|---------|
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | | 2.5 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 | 7.5 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 | 7.5 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 | 10 |

Phosphorus Flame-Retardant (PFR) Standards

| Catalog No. | Compound | Concentration | Amount |
|-------------------------|---|---------------------------|--------|
| <i>NEW</i> ULM-9693-1.2 | Resorcinol bis(diphenyl phosphate) (unlabeled) | 100 µg/mL in toluene | 1.2 mL |
| <i>NEW</i> DLM-8074-1.2 | Triethyl phosphate (D ₁₅ , 98%) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9032-1.2 | Triethyl phosphate (unlabeled) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-8901-1.2 | Tripropyl phosphate (D ₂₁ , 98%) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9090-1.2 | Tripropyl phosphate (unlabeled) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-3940-1.2 | Tributyl phosphate (D ₂₇ , 98%) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9033-1.2 | Tributyl phosphate (unlabeled) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-9070-1.2 | Triphenyl phosphate (D ₁₅ , 98%) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9091-1.2 | Triphenyl phosphate (unlabeled) | 1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-9313-1.2 | Tris(2-chloroethyl) phosphate (D ₁₂ , 98%) | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9314-1.2 | Tris(2-chloroethyl) phosphate (unlabeled) | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-9317-1.2 | Tris(2-chloroisopropyl) phosphate (D ₁₈ , 98%) | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9318-1.2 | Tris(2-chloroisopropyl) phosphate (unlabeled) | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> DLM-9315-1.2 | Tris(1,3-dichloro-2-propyl) phosphate (D ₁₅ , 98%) | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9316-1.2 | Tris(1,3-dichloro-2-propyl) phosphate (unlabeled) | 100 µg/mL in acetonitrile | 1.2 mL |

Phosphorus Flame-Retardant PFR Standard Mixtures

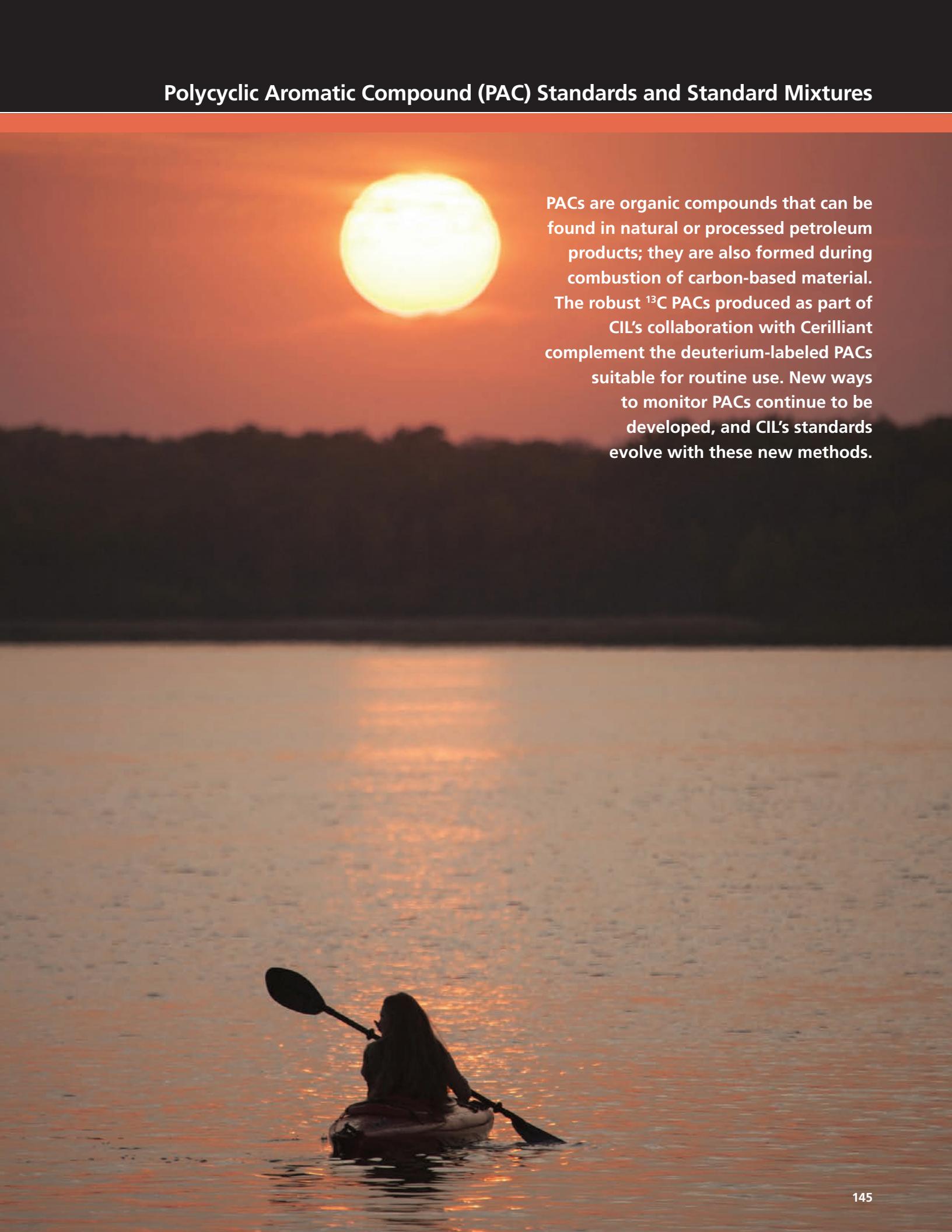
| | | |
|--------------------|--|------------------------|
| <i>NEW</i> ES-5529 | Phosphorus Flame-Retardant Standard Mixture (D, 98%) | 1.2 mL in acetonitrile |
|--------------------|--|------------------------|

| Labeled | (µg/mL) |
|---|---------|
| Tris(2-chloroethyl) phosphate (D ₁₂ , 98%) | 10 |
| Tris(2-chloroisopropyl) phosphate (D ₁₈ , 98%) | 10 |
| Tris(1,3-dichloro-2-propyl) phosphate (D ₁₅ , 98%) | 10 |
| Triethyl phosphate (D ₁₅ , 98%) | 10 |
| Tripropyl phosphate (D ₂₁ , 98%) | 10 |
| Tributyl phosphate (D ₂₇ , 98%) | 10 |
| Triphenyl phosphate (D ₁₅ , 98%) | 10 |

| | | |
|--------------------|---|------------------------|
| <i>NEW</i> ES-5530 | Phosphorus Flame-Retardant Standard Mixture | 1.2 mL in acetonitrile |
|--------------------|---|------------------------|

| Unlabeled | (µg/mL) |
|---------------------------------------|---------|
| Tris(2-chloroethyl) phosphate | 10 |
| Tris(2-chloroisopropyl) phosphate | 10 |
| Tris(1,3-dichloro-2-propyl) phosphate | 10 |
| Triethyl phosphate | 10 |
| Tripropyl phosphate | 10 |
| Tributyl phosphate | 10 |
| Triphenyl phosphate | 10 |

Polycyclic Aromatic Compound (PAC) Standards and Standard Mixtures

A photograph of a person in a kayak on a calm body of water during sunset. The sun is a large, bright orange and yellow orb in the upper left sky, casting a warm glow over the water. The horizon line is visible in the distance, and the overall scene is peaceful and scenic.

PACs are organic compounds that can be found in natural or processed petroleum products; they are also formed during combustion of carbon-based material. The robust ^{13}C PACs produced as part of CIL's collaboration with Cerilliant complement the deuterium-labeled PACs suitable for routine use. New ways to monitor PACs continue to be developed, and CIL's standards evolve with these new methods.

¹³C-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standards

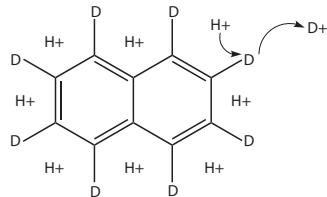
CIL, in cooperation with Cerilliant Corporation, is pleased to offer ¹³C-labeled polycyclic aromatic hydrocarbons (PAHs) as a superior alternative to deuterated standards. Although CIL has traditionally produced high-quality deuterated PAH analogs, some analysts have observed back-exchange of proton for deuterium under harsh extraction conditions and in certain matrices. If precise quantitation is required, or complete recovery information is needed, the non-exchangeable ¹³C isotope label is the right standard to use.

Deuterium Back-Exchange

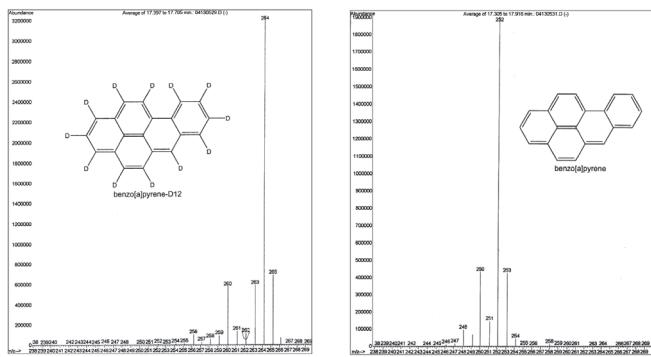
While analysts have been using deuterated PAH standards for years, labile deuterons are susceptible to back-exchange. The phenomenon is particularly likely to occur in acidic or catalytic matrices, when the importance of a reliable internal standard is greatest.

Deuterium-labeled PAH metabolites are even more susceptible to the phenomena of back-exchange and loss of protons/deuterons in the mass spectrometer.

Naphthalene-D₈ deuterium-exchange



Deuterated PAH mass spectra differ from unlabeled mass spectra



Similar Mass Spectra

Even at lower voltages the parent ion loses a considerable number of protons or deuterons. Mass spectra of ¹³C PAHs will show a succession of proton losses (M-1, M-2, M-3, M-4 etc., similar to native PAHs), while mass spectra of deuterated PAHs will show a succession of deuteron losses (M-2, M-4, M-6, M-8, etc.).

In the chromatogram for the deuterated benzo[a]pyrene, the proton losses at M-2, M-4, etc. are supplemented with proton losses of M-1, M-3, etc. This represents a loss of deuterons from incompletely deuterated species. As a result, the profile of the deuterated material does not correspond exactly to that of the unlabeled material. ¹³C-labeled benzo[a]pyrene, however, will match the unlabeled material with the 4 AMU shift being the only difference between the two profiles.

Hydroxy PAHs

PAH exposure occurs through ingestion, inhalation, and dermal contact. In the body, these compounds are predominantly metabolized as epoxides, which are converted to phenol (hydroxy) and dihydrodiol derivatives. The hydroxylated metabolites of the PAHs are excreted in human urine both as free hydroxylated metabolites and as hydroxylated metabolites conjugated to glucuronic acid or sulfate. During 1999-2002, the Centers for Disease Control and Prevention (CDC) began to measure 22 hydroxylated PAH urinary metabolites as part of the Third National Report on Human Exposure to Environmental Chemicals in the National Health and Nutrition Examination Survey (NHANES). Since then, CDC and others have focused human exposure studies on 10 target hydroxylated PAH urinary metabolites. CIL has produced isotopically labeled and unlabeled standard solutions and standard mixtures for each of these target metabolites.

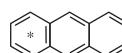
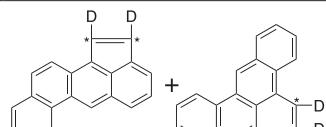
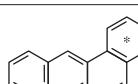
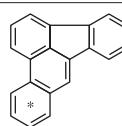
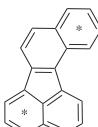
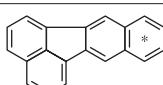
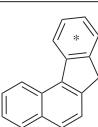
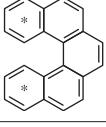
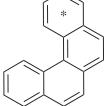
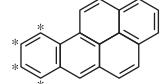
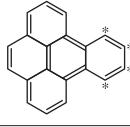
Substituted PAHs

For many years, researchers have studied the effects of diesel combustion engines, and one of the major issues related to this is the production of nitrated PAHs. Nitrated PAHs with >3 rings have been linked to the mutagenic properties associated with diesel-exhaust particulate matter. Alkyl PAHs are found in petroleum products and are typically studied in environmental forensics and fingerprinting. Alkyl PAHs may be found in high levels compared to parent PAHs and are often more persistent and bioaccumulative. Halogenated PAHs are likely the least studied but hold similar interest to parent PAHs and other substituted PAHs. Halogenated PAHs are often associated with incomplete incineration of municipal, industrial, and electronic wastes.

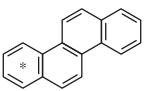
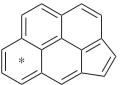
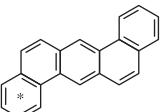
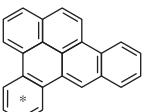
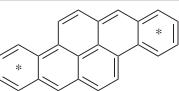
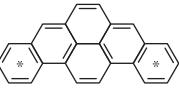
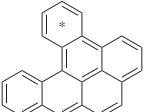
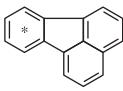
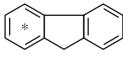
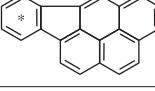
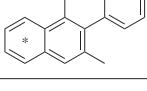
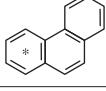
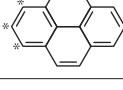
Benzo[a]Pyrene Tetrol Metabolites

Similarly to hydroxy PAHs, researchers have studied benzo[a]pyrene-7,8,9,10-tetrahydrotetrols levels in blood samples as markers for human exposure to PAHs, in particular, exposure to cigarette smoke or occupational exposure to benzo[a]pyrene, one of the most commonly studied PAHs.

¹³C-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|--------------------------|---|--|--------------------------|--------|
| CLM-1643-1.2 | Acenaphthene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-2477-1.2 | Acenaphthylene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-1333-1.2 | Anthracene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CDLM-9731-1.2 | Benz[e]aceanthrylene/Benz[j]aceanthrylene (¹³ C ₂ , 94%; D ₂ , 94%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3602-1.2 | Benz[a]anthracene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3599-1.2 | Benzo[b]fluoranthene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9590-1.2 | Benzo[j]fluoranthene (¹³ C ₁₂ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3756-1.2 | Benzo[k]fluoranthene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9363-1.2 | Benzo[c]fluorene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9610-1.2 | Benzo[b]furan (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-1364-1.2 | Benzo[ghi]perylene (¹³ C ₁₂ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9730-1.2 | Benzo[c]phenanthrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-2722-1.2 | Benzo[a]pyrene (¹³ C ₄ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-6170-1.2 | Benzo[e]pyrene (¹³ C ₄ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |

¹³C-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Structure | Concentration | Amount |
|---|--|--|---|--------|
| CLM-3757-1.2 | Chrysene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9167-1.2 | Cyclopenta[cd]pyrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3598-1.2 | Dibenz[a,h]anthracene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3835-1.2 | Dibenzo[a,e]pyrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9165-T-1.2 | Dibenzo[a,h]pyrene (¹³ C ₁₂ , 99%) |  | 100 ± 10 µg/mL in toluene | 1.2 mL |
| NEW CLM-3774-A CLM-3774-A-T-1.2 | Dibenzo[a,i]pyrene (¹³ C ₁₂ , 99%) |  | 50 ± 5 µg/mL in nonane 100 ± 10 µg/mL in toluene | 1.2 mL |
| NEW CLM-9499-1.2 | Dibenzo[a,l]pyrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3597-1.2 | Fluoranthene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3596-1.2 | Fluorene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3600-1.2 | Indeno[1,2,3-cd]pyrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| NEW CLM-9729-1.2 | 5-Methylchrysene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-1332-1.2 | Naphthalene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-2451-1.2 | Phenanthrene (¹³ C ₆ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |
| CLM-3601-1.2 | Pyrene (¹³ C ₃ , 99%) |  | 100 ± 10 µg/mL in nonane | 1.2 mL |

Deuterium-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------|--|------------------------------------|---|------------------|
| DLM-108-1.2 | Acenaphthene (D ₁₀ , 98%) | C ₁₂ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-108-0.1 | | | neat | 1 g |
| DLM-108-1 | | | neat | 5 g |
| DLM-108-5 | | | | |
| DLM-2204-1.2 | Acenaphthylene (D ₈ , 98%) | C ₁₂ D ₈ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-2204-0.1 | | | | |
| DLM-849-0.1 | Acridine (D ₉ , 98%) | C ₁₃ D ₉ N | neat | 0.1 g |
| DLM-849-0.5 | | | neat | 0.5 g |
| DLM-102-1.2 | Anthracene (D ₁₀ , 98%) | C ₁₄ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 1 g |
| DLM-102-1 | | | neat | 5 g |
| DLM-102-5 | | | | |
| DLM-610-1.2 | Benz[a]anthracene (D ₁₂ , 98%) | C ₁₈ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-610-0.1 | | | | |
| DLM-2136-1.2 | Benzo[b]fluoranthene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.01 g |
| DLM-2136-0.01 | | | | |
| DLM-1923-1.2 | Benzo[k]fluoranthene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.01 g |
| DLM-1923-0.01 | | | | |
| DLM-2135-1.2 | Benzo[ghi]perylene (D ₁₂ , 98%) | C ₂₂ D ₁₂ | 200 µg/mL in toluene-D ₈ neat | 1.2 mL 0.01 g |
| DLM-2135-0.01 | | | | |
| DLM-258-1.2 | Benzo[a]pyrene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.01 g |
| DLM-258-0.01 | | | neat | 0.05 g |
| DLM-258-0.05 | | | neat | 0.1 g |
| DLM-258-0.1 | | | | |
| DLM-257-1.2 | Benzo[e]pyrene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.01 g |
| DLM-257-0.01 | | | | |
| DLM-261-1.2 | Chrysene (D ₁₂ , 98%) | C ₁₈ D ₁₂ | 200 µg/mL in toluene-D ₈ neat | 1.2 mL 0.1 g |
| DLM-261-0.1 | | | neat | 1 g |
| DLM-261-1 | | | | |
| DLM-2715-1.2 | Coronene (D ₁₂ , 97%) | C ₂₄ D ₁₂ | 200 µg/mL in benzene neat | 1.2 mL 0.01 g |
| DLM-2715-0.01 | | | neat | 0.1 g |
| DLM-2715-0.1 | | | | |
| DLM-3843-1.2 | Dibenz[a,J]acridine (D ₁₃ , 98%) | C ₂₁ D ₁₃ N | 50 µg/mL in toluene-D ₈ | 1.2 mL |
| NEW DLM-8020-1.2 | Dibenz[a,c]anthracene (D ₁₄ , 98%) | C ₂₂ D ₁₄ | 200 µg/mL in toluene-D ₈ | 1.2 mL |
| DLM-677-1.2 | Dibenz[a,h]anthracene (D ₁₄ , 98%) | C ₂₂ D ₁₄ | 200 µg/mL in toluene-D ₈ neat | 1.2 mL 0.1 g |
| DLM-677-0.1 | | | | |
| DLM-3740-1.2 | Dibenzo[a,i]pyrene (D ₁₄ , 98%) | C ₂₄ D ₁₄ | 200 µg/mL in toluene-D ₈ | 1.2 mL |
| DLM-3841-1.2 | 7H-Dibenzo[c,g]carbazole (D ₁₂ , 98%) | C ₂₀ D ₁₂ NN | 50 µg/mL in toluene-D ₈ | 1.2 mL |
| DLM-2140-1.2 | Fluoranthene (D ₁₀ , 98%) | C ₁₆ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-2140-0.1 | | | | |
| DLM-1123-1.2 | Fluorene (D ₁₀ , 98%) | C ₁₃ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-1123-0.1 | | | neat | 1 g |
| DLM-1123-1 | | | | |
| DLM-2148-1.2 | Indeno[1,2,3-cd]pyrene (D ₁₂ , 98%) | C ₂₂ D ₁₂ | 200 µg/mL in isoctane neat | 1.2 mL 0.01 g |
| DLM-2148-0.01 | | | | |
| DLM-365-1.2 | Naphthalene (D ₈ , 99%) | C ₁₀ D ₈ | 200 µg/mL in isoctane neat | 1.2 mL 1 g |
| DLM-365-1 | | | neat | 5 g |
| DLM-365-5 | | | neat | 10 g |
| DLM-365-10 | | | neat | |
| DLM-3875-10 | Naphthalene (D ₈ , 99.5%) | C ₁₀ D ₈ | neat | 10 g |
| DLM-366-1.2 | Perylene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in toluene-D ₈ neat | 1.2 mL 0.1 g |
| DLM-366-0.1 | | | neat | 1 g |
| DLM-366-1 | | | | |
| DLM-371-1.2 | Phenanthrene (D ₁₀ , 98%) | C ₁₄ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-371-0.1 | | | neat | 1 g |
| DLM-371-1 | | | neat | 5 g |
| DLM-371-5 | | | neat | |
| DLM-155-1.2 | Pyrene (D ₁₀ , 98%) | C ₁₆ D ₁₀ | 200 µg/mL in isoctane neat | 1.2 mL 0.1 g |
| DLM-155-0.1 | | | neat | 0.5 g |
| DLM-155-0.5 | | | | |

Deuterium-Labeled PAH Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------------|-------------------------------------|---|-----------------------|--------|
| NEW DLM-450-1 | o-Terphenyl (D ₁₄ , 98%) | C ₆ D ₄ (C ₆ D ₅) ₂ | neat | 1 g |
| | DLM-450-5 | | neat | 5 g |
| NEW DLM-382-1.2 | p-Terphenyl (D ₁₄ , 98%) | C ₆ D ₄ (C ₆ D ₅) ₂ | 200 µg/mL in isoctane | 1.2 mL |
| | DLM-382-1 | | neat | 1 g |
| NEW DLM-382-5 | DLM-382-5 | | neat | 5 g |
| | DLM-601-0.1 | C ₁₈ D ₁₂ | neat | 0.1 g |
| DLM-601-1 | | | neat | 1 g |

Unlabeled PAH Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------------|---|-----------------------------------|-----------------------------------|-----------------------|
| ULM-7413-1.2 | Acenaphthene | C ₁₂ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7422-1.2 | Acenaphthylene | C ₁₂ H ₈ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7412-1.2 | Anthracene | C ₁₄ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL |
| NEW ULM-9758-1.2 | Benz[e]aceanthrylene/Benz[j]aceanthrylene | C ₂₀ H ₁₂ | 200 µg/mL in nonane | 1.2 mL |
| | ULM-2415-I-1.2 | Benz[a]anthracene | C ₁₈ H ₁₂ | 200 µg/mL in isoctane |
| ULM-2415-0.1 | | | neat | 0.1 g |
| ULM-2416-I-1.2 | Benzo[b]fluoranthene | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-2416-0.1 | | | neat | 0.1 g |
| NEW ULM-2411-1.2 | Benzo[j]fluoranthene | C ₂₀ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-2411-25 | | neat | 25 mg |
| NEW ULM-2417-I-1.2 | Benzo[k]fluoranthene | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | ULM-2417-0.1 | | neat | 0.1 g |
| NEW ULM-9357-1.2 | Benzo[c]fluorene | C ₁₇ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9618-1.2 | Benzo[b]furan | C ₈ H ₆ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2418-1.2 | Benzo[ghi]perylene | C ₂₂ H ₁₂ | 200 µg/mL in toluene | 1.2 mL |
| ULM-2418-0.1 | | | neat | 0.1 g |
| NEW ULM-8155-1.2 | Benzo[c]phenanthrene | C ₁₈ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-8155-25 | | neat | 25 mg |
| ULM-2412-I-1.2 | Benzo[a]pyrene | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-2412-0.1 | | | neat | 0.1 g |
| ULM-7423-1.2 | Benzo[e]pyrene | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7424-1.2 | Chrysene | C ₁₈ H ₁₂ | 200 µg/mL in toluene | 1.2 mL |
| ULM-6576-1.2 | Coronene | C ₂₄ H ₁₂ | 200 µg/mL in benzene | 1.2 mL |
| NEW ULM-6891-1.2 | Cyclopenta[cd]pyrene | C ₁₈ H ₁₀ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-3884-1.2 | Dibenz[a,j]acridine | C ₂₁ H ₁₃ N | 50 µg/mL in toluene |
| ULM-3884-25 | | | neat | 25 mg |
| NEW ULM-9759-1.2 | Dibenz[a,c]anthracene | C ₂₂ H ₁₄ | 200 µg/mL in toluene | 1.2 mL |
| | ULM-2422-T-1.2 | Dibenz[a,h]anthracene | C ₂₂ H ₁₄ | 200 µg/mL in toluene |
| ULM-2422-0.1 | | | neat | 0.1 g |
| ULM-3885-1.2 | 7H-Dibenzo[c,g]carbazole | C ₂₀ H ₁₃ N | 50 µg/mL in toluene | 1.2 mL |
| ULM-6671-1.2 | Dibenzo[a,e]fluoranthene | C ₂₄ H ₁₄ | 200 µg/mL in toluene | 1.2 mL |
| NEW ULM-1226-1.2 | Dibenzo[a,e]pyrene | C ₂₄ H ₁₄ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-1226-0.01 | | neat | 0.01 g |
| ULM-1227-T-1.2 | Dibenzo[a,h]pyrene | C ₂₄ H ₁₄ | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-1227-0.01 | | | neat | 0.01 g |
| ULM-2423-1.2 | Dibenzo[a,i]pyrene | C ₂₄ H ₁₄ | 200 µg/mL in toluene | 1.2 mL |
| NEW ULM-2423-A-1.2 | | | 50 µg/mL in nonane | 1.2 mL |
| ULM-1253-1.2 | Dibenzo[a,l]pyrene | C ₂₄ H ₁₄ | 200 µg/mL in toluene | 1.2 mL |
| ULM-1253-25 | | | neat | 25 mg |
| ULM-7421-1.2 | Fluoranthene | C ₁₆ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7414-1.2 | Fluorene | C ₁₃ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-2426-I-1.2 | Indeno[1,2,3-cd]pyrene | C ₂₂ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-2426-25 | | | neat | 25 mg |
| ULM-7425-1.2 | Naphthalene | C ₁₀ H ₈ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7426-1.2 | Perylene | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7427-1.2 | Phenanthrene | C ₁₄ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7417-1.2 | Pyrene | C ₁₆ H ₁₀ | 200 µg/mL in toluene | 1.2 mL |
| ULM-7428-1.2 | p-Terphenyl | C ₁₈ H ₁₄ | 200 µg/mL in isoctane | 1.2 mL |

¹³C-Labeled PAH Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|------------------|
| ES-4087 | EPA 16 PAH Cocktail | 1.2 mL in nonane |
| | Labeled | (μ g/mL) |
| | Acenaphthene (¹³ C ₆ , 99%) | 5 |
| | Acenaphthylene (¹³ C ₆ , 99%) | 5 |
| | Anthracene (¹³ C ₆ , 99%) | 5 |
| | Benz[a]anthracene (¹³ C ₆ , 99%) | 5 |
| | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Benzo[k]fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Benzo[ghi]perylene (¹³ C ₁₂ , 99%) | 5 |
| | Benzo[a]pyrene (¹³ C ₄ , 99%) | 5 |
| | Chrysene (¹³ C ₆ , 99%) | 5 |
| | Dibenz[a,h]anthracene (¹³ C ₆ , 99%) | 5 |
| | Fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Fluorene (¹³ C ₆ , 99%) | 5 |
| | Indeno[1,2,3-cd]pyrene (¹³ C ₆ , 99%) | 5 |
| | Naphthalene (¹³ C ₆ , 99%) | 5 |
| | Phenanthrene (¹³ C ₆ , 99%) | 5 |
| | Pyrene (¹³ C ₃ , 99%) | 5 |
| NEW | ES-5546 Custom 16 PAH Solution | 1.2 mL in nonane |
| | Labeled | (μ g/mL) |
| | Naphthalene (¹³ C ₁₀ , 99%) | 10 |
| | Benzo[c]phenanthrene (¹³ C ₆ , 99%) | 1 |
| | Cyclopenta[cd]pyrene (¹³ C ₆ , 99%) | 1 |
| | Benz[a]anthracene (¹³ C ₆ , 99%) | 1 |
| | Chrysene (¹³ C ₆ , 99%) | 1 |
| | 5-Methylchrysene (¹³ C ₆ , 99%) | 1 |
| | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | 1 |
| | Benzo[k]fluoranthene (¹³ C ₆ , 99%) | 1 |
| | Benz[e]aceanthrylene/Benz[j]aceanthrylene (¹³ C ₂ , 94%; D ₂ , 94%) | 1 |
| | Benzo[a]pyrene (¹³ C ₄ , 99%) | 1 |
| | Indeno[1,2,3-cd]pyrene (¹³ C ₆ , 99%) | 1 |
| | Dibenz[a,h]anthracene (¹³ C ₆ , 99%) | 1 |
| | Dibenzo[a,l]pyrene (¹³ C ₆ , 99%) | 1 |
| | Dibenzo[a,e]pyrene (¹³ C ₆ , 99%) | 1 |
| | Dibenzo[a,i]pyrene (¹³ C ₁₂ , 99%) | 1 |
| | Dibenzo[a,h]pyrene (¹³ C ₁₂ , 99%) | 1 |
| NEW | ES-5539 EFSA-8 ¹³ C PAH Standard Mixture | 1.2 mL in nonane |
| | Labeled | (μ g/mL) |
| | Benz[a]anthracene (¹³ C ₆ , 99%) | 5 |
| | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Benzo[k]fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Benzo[ghi]perylene (¹³ C ₁₂ , 99%) | 5 |
| | Benzo[a]pyrene (¹³ C ₄ , 99%) | 5 |
| | Chrysene (¹³ C ₆ , 99%) | 5 |
| | Dibenz[a,h]anthracene (¹³ C ₆ , 99%) | 5 |
| | Indeno[1,2,3-cd]pyrene (¹³ C ₆ , 99%) | 5 |
| NEW | ES-5540 EFSA-4 ¹³ C PAH Standard Mixture | 1.2 mL in nonane |
| | Labeled | (μ g/mL) |
| | Benz[a]anthracene (¹³ C ₆ , 99%) | 5 |
| | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | 5 |
| | Benzo[a]pyrene (¹³ C ₄ , 99%) | 5 |
| | Chrysene (¹³ C ₆ , 99%) | 5 |

Deuterium-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|-----------------------|--|--|
| ES-2528 | PAH Cocktail for CARB Method 429 | 1 mL in benzene-D ₆ |
| | Labeled | (μ g/mL) |
| | Acenaphthene (D ₁₀ , 98%) | 100 |
| | Acenaphthylene (D ₈ , 98%) | 100 |
| | Anthracene (D ₁₀ , 98%) | 100 |
| | Benz[a]anthracene (D ₁₂ , 98%) | 100 |
| | Benzo[b]fluoranthene (D ₁₂ , 98%) | 100 |
| | Benzo[k]fluoranthene (D ₁₂ , 98%) | 100 |
| | Benzo[ghi]perylene (D ₁₂ , 98%) | 100 |
| | Benzo[a]pyrene (D ₁₂ , 98%) | 100 |
| | Chrysene (D ₁₂ , 98%) | 100 |
| | Dibenz[a,h]anthracene (D ₁₄ , 98%) | 100 |
| | Fluoranthene (D ₁₀ , 98%) | 100 |
| | Fluorene (D ₁₀ , 98%) | 100 |
| | Indeno[1,2,3-cd]pyrene (D ₁₂ , 98%) | 100 |
| | Naphthalene (D ₈ , 99%) | 100 |
| | Phenanthrene (D ₁₀ , 98%) | 100 |
| | Pyrene (D ₁₀ , 98%) | 100 |
| ES-5164 NEW | PAH Surrogate Standard Mixture | 10 mL in 90% toluene/10% isoctane |
| ES-5164-1.2 | | 1.2 mL in 90% toluene/10% isoctane |
| | Labeled | (μ g/mL) |
| | Naphthalene (D ₈ , 99%) | 200 |
| | Benz[a]anthracene (D ₁₂ , 98%) | 200 |
| | Phenanthrene (D ₁₀ , 98%) | 200 |
| | Fluoranthene (D ₁₀ , 98%) | 200 |
| | Benzo[b]fluoranthene (D ₁₂ , 98%) | 200 |
| | Benzo[a]pyrene (D ₁₂ , 98%) | 200 |
| | Benzo[ghi]perylene (D ₁₂ , 98%) | 200 |
| | Indeno[1,2,3-cd]pyrene (D ₁₂ , 98%) | 200 |
| | Dibenz[a,h]anthracene (D ₁₄ , 98%) | 200 |
| | Acenaphthylene (D ₈ , 98%) | 200 |
| | Acenaphthene (D ₁₀ , 98%) | 200 |
| | Fluorene (D ₁₀ , 98%) | 200 |
| | Pyrene (D ₁₀ , 98%) | 200 |
| | Benzo[k]fluoranthene (D ₁₂ , 98%) | 200 |
| | Perylene (D ₁₂ , 98%) | 200 |
| | Chrysene (D ₁₂ , 98%) | 200 |
| ES-2044 | PAH Surrogate Cocktail | 1 mL in 50% MeCl-D ₂ /50% methanol-OD |
| | Labeled | (μ g/mL) |
| | Acenaphthylene (D ₈ , 98%) | 200 |
| | Benzo[a]pyrene (D ₁₂ , 98%) | 200 |
| | Benzo[ghi]perylene (D ₁₂ , 98%) | 200 |
| | Fluoranthene (D ₁₀ , 98%) | 200 |
| | Naphthalene (D ₈ , 99%) | 200 |
| | Phenanthrene (D ₁₀ , 98%) | 200 |
| | Pyrene (D ₁₀ , 98%) | 200 |

Deuterium-Labeled Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|--|-----------------------------------|--|
| NEW ES-5481 | PAH Mixture | 5 mL in toluene |
| Labeled (µg/mL) | | |
| Chrysene (D ₁₂ , 98%) | 2500 | |
| Dibenz[a,h]anthracene (D ₁₄ , 98%) | 2500 | |
| Naphthalene (D ₈ , 99%) | 2500 | |
| Perylene (D ₁₂ , 98%) | 2500 | |
| Phenanthrene (D ₁₀ , 98%) | 2500 | |
| Acenaphthene (D ₁₀ , 98%) | 2500 | |
| <hr/> | | |
| ES-2043 | "EEC Six" PAH Cocktail | 1.2 mL in benzene-D ₆ |
| Labeled (µg/mL) | | |
| Benzo[b]fluoranthene (D ₁₂ , 98%) | 1000 | |
| Benzo[k]fluoranthene (D ₁₂ , 98%) | 1000 | |
| Benzo[ghi]perylene (D ₁₂ , 98%) | 1000 | |
| Benzo[a]pyrene (D ₁₂ , 98%) | 1000 | |
| Indeno[1,2,3-cd]pyrene (D ₁₂ , 98%) | 1000 | |
| Fluoranthene (D ₁₀ , 98%) | 1000 | |
| <hr/> | | |
| ES-5386 | PAH-SIM Recovery Standard Mixture | 1.2 mL in MeCl-D ₂ |
| Labeled (µg/mL) | | |
| 2-Methylnaphthalene (D ₁₀ , 98%) | 1000 | |
| Anthracene (D ₁₀ , 98%) | 1000 | |
| p-Terphenyl (D ₁₄ , 98%) | 1000 | |
| Benzo[e]pyrene (D ₁₂ , 98%) | 1000 | |
| <hr/> | | |
| NEW ES-5498 | PAH Two-Component Mixture | 1.2 mL in 80% isoctane/ 20% toluene |
| Labeled (µg/mL) | | |
| Benzo[a]pyrene (D ₁₂ , 98%) | 2000 | |
| Fluoranthene (D ₁₀ , 98%) | 2000 | |
| <hr/> | | |
| NEW ES-9463 | PAH Injection Standard | 10 mL in isoctane |
| Labeled (ng/mL) | | |
| Anthracene (D ₁₀ , 98%) | 10 | |
| Fluoranthene (D ₁₀ , 98%) | 10 | |
| Benz[a]anthracene (D ₁₂ , 98%) | 10 | |
| <hr/> | | |
| NEW ES-9464 | PAH Recovery Standard | 10 mL in isoctane/toluene-D ₈ |
| Labeled (ng/mL) | | |
| Acenaphthene (D ₁₀ , 98%) | 10 | |
| p-Terphenyl (D ₁₄ , 98%) | 10 | |
| Perylene (D ₁₂ , 98%) | 10 | |

Unlabeled Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------------|----------------------------------|-------------------|
| NEW ES-5437 | PAH Cocktail for CARB Method 429 | 1.2 mL in benzene |

| Unlabeled | (μ g/mL) |
|------------------------|---------------|
| Acenaphthene | 100 |
| Acenaphthylene | 100 |
| Anthracene | 100 |
| Benz[a]anthracene | 100 |
| Benzo[b]fluoranthene | 100 |
| Benzo[k]fluoranthene | 100 |
| Benzo[ghi]perylene | 100 |
| Benzo[a]pyrene | 100 |
| Chrysene | 100 |
| Dibenz[a,h]anthracene | 100 |
| Fluoranthene | 100 |
| Fluorene | 100 |
| Indeno[1,2,3-cd]pyrene | 100 |
| Naphthalene | 100 |
| Phenanthrene | 100 |
| Pyrene | 100 |

| | | |
|--------------------|---------------------------------------|------------------|
| NEW ES-5549 | Custom 16 PAH Native Standard Mixture | 1.2 mL in nonane |
|--------------------|---------------------------------------|------------------|

| Unlabeled | (μ g/mL) |
|---|---------------|
| Naphthalene | 10 |
| Benzo[c]phenanthrene | 1 |
| Cyclopenta[cd]pyrene | 1 |
| Benz[a]anthracene | 1 |
| Chrysene | 1 |
| 5-Methylchrysene | 1 |
| Benzo[b]fluoranthene | 1 |
| Benzo[k]fluoranthene | 1 |
| Benz[e]aceanthrylene/Benz[j]aceanthrylene | 1 |
| Benzo[a]pyrene | 1 |
| Indeno[1,2,3-cd]pyrene | 1 |
| Dibenz[a,h]anthracene | 1 |
| Dibenzo[a,l]pyrene | 1 |
| Dibenzo[a,e]pyrene | 1 |
| Dibenzo[a,i]pyrene | 1 |
| Dibenzo[a,h]pyrene | 1 |

| | | |
|--------------------|------------------------------------|--------------------|
| NEW ES-5541 | EFSA-8 Native PAH Standard Mixture | 1.2 mL in isoctane |
|--------------------|------------------------------------|--------------------|

| Unlabeled | (μ g/mL) |
|------------------------|---------------|
| Benz[a]anthracene | 5 |
| Benzo[b]fluoranthene | 5 |
| Benzo[k]fluoranthene | 5 |
| Benzo[ghi]perylene | 5 |
| Benzo[a]pyrene | 5 |
| Chrysene | 5 |
| Dibenz[a,h]anthracene | 5 |
| Indeno[1,2,3-cd]pyrene | 5 |

Unlabeled Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|---------------------------------------|------------------------------------|------------------------------------|
| NEW ES-5542 | EFSA-4 Native PAH Standard Mixture | 1.2 mL in isoctane |
| Unlabeled ($\mu\text{g/mL}$) | | |
| Benz[a]anthracene | 5 | |
| Benzo[b]fluoranthene | 5 | |
| Benzo[a]pyrene | 5 | |
| Chrysene | 5 | |
| <hr/> | | |
| ES-5438 | PAH Native Standard Mixture | 1.2 mL in 90% toluene/10% isoctane |
| Unlabeled ($\mu\text{g/mL}$) | | |
| Naphthalene | 200 | |
| Benz[a]anthracene | 200 | |
| Phenanthrene | 200 | |
| Fluoranthene | 200 | |
| Benzo[b]fluoranthene | 200 | |
| Benzo[a]pyrene | 200 | |
| Benzo[ghi]perylene | 200 | |
| Indeno[1,2,3- <i>cd</i>]pyrene | 200 | |
| Dibenz[a,h]anthracene | 200 | |
| Acenaphthylene | 200 | |
| Acenaphthene | 200 | |
| Fluorene | 200 | |
| Pyrene | 200 | |
| Benzo[k]fluoranthene | 200 | |
| Perylene | 200 | |
| Chrysene | 200 | |
| <hr/> | | |
| NEW ES-5503 | PAH-SIM Recovery Standard Mixture | 1.2 mL in MeCl |
| Unlabeled ($\mu\text{g/mL}$) | | |
| 2-Methylnaphthalene | 1000 | |
| Anthracene | 1000 | |
| <i>p</i> -Terphenyl | 1000 | |
| Benzo[e]pyrene | 1000 | |

Alkyl Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--------------------|----------------------------|--------|
| DLM-2845-1.2 | 9,10-Dimethylanthracene (D_{14} , 98%) | $C_{16}D_{14}$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-6234-1.2 | 9,10-Dimethylanthracene (unlabeled) | $C_{16}H_{14}$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-2852-1.2 | 1,6-Dimethylnaphthalene (D_{12} , 98%) | $C_{12}D_{12}$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-6182-1.2 | 1,6-Dimethylnaphthalene (unlabeled) | $C_{12}H_{12}$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-2854-1.2 | 1,8-Dimethylnaphthalene (D_{12} , 98%) | $C_{12}D_{12}$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| ULM-6181-1.2 | 1,8-Dimethylnaphthalene (unlabeled) | $C_{12}H_{12}$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-2853-1.2 | 2,6-Dimethylnaphthalene (D_{12} , 98%) | $C_{12}D_{12}$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| ULM-7271-1.2 | 2,6-Dimethylnaphthalene (unlabeled) | $C_{12}H_{12}$ | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-9729-1.2 | 5-Methylchrysene ($^{13}C_6$, 99%) | $*C_6C_{13}H_{14}$ | 100 µg/mL in nonane | 1.2 mL |
| DLM-3842-1.2 | 5-Methylchrysene (methyl- D_3 , 98%) | $C_{19}D_3H_{11}$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| ULM-6235-1.2 | 5-Methylchrysene (unlabeled) | $C_{19}H_{14}$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-1607-1 | 1-Methylnaphthalene (D_{10} , 98%) | $C_{11}D_{10}$ | neat | 1 g |
| CLM-3621-1.2 | 2-Methylnaphthalene ($^{13}C_6$, 99%) | $*C_6C_5H_{10}$ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1322-1.2 | 2-Methylnaphthalene (D_{10} , 98%) | $C_{11}D_{10}$ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7416-1.2 | 2-Methylnaphthalene (unlabeled) | $C_{11}H_{10}$ | 200 µg/mL in isoctane | 1.2 mL |
| NEW CLM-8174-1.2 | 2-Methyl-1-naphthol (5,6,7,8,9,10- $^{13}C_6$, 99%) | $*C_6C_5H_{10}$ | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8239-1.2 | 2-Methyl-1-naphthol (unlabeled) | $C_{11}H_{10}O$ | 50 µg/mL in toluene | 1.2 mL |

Halogenated Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|---|------------------------|---------------------|--------|
| NEW CLM-9028-1.2 | 7-Bromobenz[a]anthracene ($^{13}C_6$, 99%) | $*C_6C_{12}H_{11}Br$ | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-9025-1.2 | 7-Bromobenz[a]anthracene (unlabeled) | $C_{18}H_{11}Br$ | 50 µg/mL in toluene | 1.2 mL |
| ULM-8269-1.2 | 9-Chloroanthracene (unlabeled) | $C_{14}H_9Cl$ | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-8989-1.2 | 7-Chlorobenz[a]anthracene ($^{13}C_6$, 99%) | $*C_6C_{12}H_1Cl$ | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-9011-1.2 | 7-Chlorobenz[a]anthracene (unlabeled) | $C_{18}H_{11}Cl$ | 50 µg/mL in toluene | 1.2 mL |
| ULM-8270-1.2 | 9-Chlorophenanthrene (unlabeled) | $C_{14}H_9Cl$ | 50 µg/mL in toluene | 1.2 mL |
| CLM-8267-1.2 | 1-Chloropyrene (mix of ring labeling) ($^{13}C_6$, 99%) | $C_{10}*C_6H_9Cl$ | 50 µg/mL in toluene | 1.2 mL |
| ULM-8268-1.2 | 1-Chloropyrene (unlabeled) | $C_{16}H_9Cl$ | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-9029-1.2 | 7,12-Dichlorobenz[a]anthracene ($^{13}C_6$, 99%) | $C_{12}*C_6H_{10}Cl_2$ | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-9024-1.2 | 7,12-Dichlorobenz[a]anthracene (unlabeled) | $C_{18}H_{10}Cl_2$ | 50 µg/mL in toluene | 1.2 mL |

Nitro Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|---|--------------------|----------------------------|--------|
| DLM-3836-1.2 | 5-Nitroacenaphthene (D_9 , 98%) | $C_{12}D_9NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-8790-1.2 | 5-Nitroacenaphthene (unlabeled) | $C_{12}H_9NO_2$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-4712-1.2 | 9-Nitroanthracene (D_9 , 98%) | $C_{14}D_9NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-8365-1.2 | 9-Nitroanthracene (unlabeled) | $C_{14}H_9NO_2$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-3839-1.2 | 6-Nitrochrysene (D_{11} , 98%) | $C_{18}D_{11}NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-3881-1.2 | 6-Nitrochrysene (unlabeled) | $C_{18}H_{11}NO_2$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-4711-1.2 | 3-Nitrofluoranthene (D_9 , 98%) CP 87% | $C_{16}D_9NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| NEW ULM-6600-1.2 | 3-Nitrofluoranthene (unlabeled) CP 87% | $C_{16}H_9NO_2$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-3837-1.2 | 2-Nitrofluorene (D_9 , 98%) | $C_{13}D_9NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| ULM-3883-1.2 | 2-Nitrofluorene (unlabeled) | $C_{13}H_9NO_2$ | 50 µg/mL in toluene | 1.2 mL |
| DLM-1528-1.2 | 1-Nitropyrene (D_9 , 98%) | $C_{18}D_9NO_2$ | 50 µg/mL in toluene- D_8 | 1.2 mL |
| ULM-3978-1.2 | 1-Nitropyrene (unlabeled) | $C_{18}H_9NO_2$ | 50 µg/mL in toluene | 1.2 mL |

Benzo[a]pyrene Metabolites

| Catalog No. | Compound | Concentration | Amount |
|-------------------------|---|-----------------------|---------|
| NEW CLM-7245-1.2 | (+/-)-Benzo[a]pyrene R-7,T-8,C-9,C-10-tetrahydrotetrol (ring- $^{13}C_6$, 99%) | 100 µg/mL in methanol | Inquire |
| NEW CLM-7246-1.2 | (+/-)-Benzo[a]pyrene R-7,T-8,T-9,C-10-tetrahydrotetrol (ring- $^{13}C_6$, 99%) | 100 µg/mL in methanol | Inquire |
| NEW CLM-7308-1.2 | (+/-)-Benzo[a]pyrene R-7,T-8,C-9,T-10-tetrahydrotetrol (ring- $^{13}C_6$, 99%) | 100 µg/mL in methanol | Inquire |
| NEW CLM-7627-1.2 | (+/-)-Benzo[a]pyrene R-7,T-8,T-9,T-10-tetrahydrotetrol (ring- $^{13}C_6$, 99%) | 100 µg/mL in methanol | Inquire |

Hydroxy Polycyclic Aromatic Hydrocarbon (PAH) Standards

| Catalog No. | Compound | Concentration | Amount |
|---------------------------|--|----------------------|--------|
| CLM-4860-T-1.2 | 6-Hydroxychrysene (mix of ring labeling) (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-7552-1.2 | 6-Hydroxychrysene (unlabeled) | 50 µg/mL in isoctane | 1.2 mL |
| NEW CLM-6087-1.2 | 2-Hydroxyfluorene (random- ¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8973-1.2 | 2-Hydroxyfluorene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-8977-1.2 | 3-Hydroxyfluorene (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8974-1.2 | 3-Hydroxyfluorene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-7700-1.2 | 9-Hydroxyfluorene (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8975-1.2 | 9-Hydroxyfluorene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-7701-1.2 | 1-Hydroxynaphthalene (1-naphthol) (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8971-1.2 | 1-Hydroxynaphthalene (1-naphthol) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-7713-1.2 | 2-Hydroxynaphthalene (2-naphthol) (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8972-1.2 | 2-Hydroxynaphthalene (2-naphthol) (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-7669-1.2 | 1-Hydroxyphenanthrene (¹³ C ₄ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-7929-1.2 | 1-Hydroxyphenanthrene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-8463-T-1.2 | 2-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8464-T-1.2 | 2-Hydroxyphenanthrene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| CLM-4859-T-1.2 | 3-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-7446-1.2 | 3-Hydroxyphenanthrene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-7670-1.2 | 4-Hydroxyphenanthrene (¹³ C ₄ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-7928-1.2 | 4-Hydroxyphenanthrene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |
| NEW CLM-9012-1.2 | 1-Hydroxypyrene (¹³ C ₆ , 99%) | 50 µg/mL in toluene | 1.2 mL |
| NEW ULM-8976-1.2 | 1-Hydroxypyrene (unlabeled) | 50 µg/mL in toluene | 1.2 mL |

Hydroxy Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|---|------------------------|
| NEW ES-5472 | CDC OH-PAH Calibration Standards [CS1-CS10] | 10 x 0.5 mL in toluene |
| NEW ES-5472-CS1 | CDC OH-PAH Calibration Standard [CS1] | 0.5 mL in toluene |
| NEW ES-5472-CS2 | CDC OH-PAH Calibration Standard [CS2] | 0.5 mL in toluene |
| NEW ES-5472-CS3 | CDC OH-PAH Calibration Standard [CS3] | 0.5 mL in toluene |
| NEW ES-5472-CS4 | CDC OH-PAH Calibration Standard [CS4] | 0.5 mL in toluene |
| NEW ES-5472-CS5 | CDC OH-PAH Calibration Standard [CS5] | 0.5 mL in toluene |
| NEW ES-5472-CS6 | CDC OH-PAH Calibration Standard [CS6] | 0.5 mL in toluene |
| NEW ES-5472-CS7 | CDC OH-PAH Calibration Standard [CS7] | 0.5 mL in toluene |
| NEW ES-5472-CS8 | CDC OH-PAH Calibration Standard [CS8] | 0.5 mL in toluene |
| NEW ES-5472-CS9 | CDC OH-PAH Calibration Standard [CS9] | 0.5 mL in toluene |
| NEW ES-5472-CS10 | CDC OH-PAH Calibration Standard [CS10] | 0.5 mL in toluene |

All concentrations are in ng/mL

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 | CS10 |
|--|-------|-----|-----|-----|-----|-----|-----|------|------|------|--------|
| 1-Hydroxynaphthalene | | 4 | 8 | 20 | 40 | 200 | 400 | 2000 | 4000 | 8000 | 16,000 |
| 2-Hydroxynaphthalene | | 4 | 8 | 20 | 40 | 200 | 400 | 2000 | 4000 | 8000 | 16,000 |
| 2-Hydroxyfluorene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 3-Hydroxyfluorene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 9-Hydroxyfluorene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 1-Hydroxyphenanthrene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 2-Hydroxyphenanthrene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 3-Hydroxyphenanthrene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 4-Hydroxyphenanthrene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| 1-Hydroxypyrene | | 1 | 2 | 5 | 10 | 50 | 100 | 500 | 1000 | — | — |
| Labeled | | | | | | | | | | | |
| 1-Hydroxynaphthalene (¹³ C ₆ , 99%) | | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 |
| 2-Hydroxynaphthalene (¹³ C ₆ , 99%) | | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 | 400 |
| 2-Hydroxyfluorene (random- ¹³ C ₆ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3-Hydroxyfluorene (¹³ C ₆ , 98%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 9-Hydroxyfluorene (¹³ C ₆ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1-Hydroxyphenanthrene (¹³ C ₄ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2-Hydroxyphenanthrene (¹³ C ₆ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3-Hydroxyphenanthrene (¹³ C ₆ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4-Hydroxyphenanthrene (¹³ C ₄ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 1-Hydroxypyrene (¹³ C ₆ , 99%) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |

Hydroxy Polycyclic Aromatic Hydrocarbon (PAH) Standard Mixtures

| Catalog No. | Compound | Amount |
|--|--|-------------------|
| NEW ES-5473-T | CDC OH-PAH Spiking Standard | 0.5 mL in toluene |
| Labeled | | |
| | | (ng/mL) |
| | 1-Hydroxynaphthalene (¹³ C ₆ , 99%) | 10,000 |
| | 2-Hydroxynaphthalene (¹³ C ₆ , 99%) | 10,000 |
| | 2-Hydroxyfluorene (random- ¹³ C ₆ , 99%) | 2500 |
| | 3-Hydroxyfluorene (¹³ C ₆ , 98%) | 2500 |
| | 9-Hydroxyfluorene (¹³ C ₆ , 99%) | 2500 |
| | 1-Hydroxyphenanthrene (¹³ C ₄ , 99%) | 2500 |
| | 2-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 2500 |
| | 3-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 2500 |
| | 4-Hydroxyphenanthrene (¹³ C ₄ , 99%) | 2500 |
| | 1-Hydroxypyrene (¹³ C ₆ , 99%) | 2500 |
| NEW ES-5474 CDC PCB Recovery Standard for OH-PAHs | | |
| | | |
| Labeled | | |
| | | IUPAC |
| | | (ng/mL) |
| | 2,4,4'-TriCB (¹³ C ₁₂ , 99%) | 28 |
| | 2,3,3',4,4'-PentaCB (¹³ C ₁₂ , 99%) | 105 |
| | 2,2',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 153 |
| | 2,3',4,4',5,5'-HexaCB (¹³ C ₁₂ , 99%) | 167 |
| NEW ES-5484 CDC OH-PAH Native PAR Standard | | |
| | | |
| Unlabeled | | |
| | | (ng/mL) |
| | 1-Hydroxynaphthalene | 4000 |
| | 2-Hydroxynaphthalene | 4000 |
| | 2-Hydroxyfluorene | 1000 |
| | 3-Hydroxyfluorene | 1000 |
| | 9-Hydroxyfluorene | 1000 |
| | 1-Hydroxyphenanthrene | 1000 |
| | 2-Hydroxyphenanthrene | 1000 |
| | 3-Hydroxyphenanthrene | 1000 |
| | 4-Hydroxyphenanthrene | 1000 |
| | 1-Hydroxypyrene | 1000 |

Isotope-Labeled Polychlorinated Naphthalene (PCN) Standards

| Catalog No. | Compound | PCN | Concentration | Amount |
|-------------|---|-----|---------------------------------|---------|
| NEW | ECN-5217 2-Monochloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 2 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| NEW | ECN-5520 1,5-Dichloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 6 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| NEW | ECN-5575 1,2,3-Trichloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 13 | 10 $\mu\text{g/mL}$ in isoctane | Inquire |
| | ECN-5240 1,2,3,4-Tetrachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 27 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5241 1,3,5,7-Tetrachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 42 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5250 1,2,3,5,7-Pentachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 52 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5260 1,2,3,4,5,7-Hexachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 64 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| NEW | ECN-5267 1,2,3,4,5,8-Hexachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 65 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5261-A 1,2,3,5,6,7-Hexachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 67 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5270-A 1,2,3,4,5,6,7-Heptachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 73 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| | ECN-5280 Octachloronaphthalene ($^{13}\text{C}_{10}$, 99%) | 75 | 10 $\mu\text{g/mL}$ in isoctane | 1.2 mL |
| NEW | DLM-2005-1.2 2-Chloronaphthalene (D ₇ , 98%) | | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW | DLM-2005-0.01 | | neat | 0.01 g |
| NEW | DLM-2005-0.1 | | neat | 0.1 g |

Unlabeled Polychlorinated Naphthalene (PCN) Standards

| Catalog No. | Compound | PCN | Concentration | Amount |
|-------------|---|-----|--------------------------------|---------|
| | ECN-2610 1-Monochloronaphthalene | 1 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2611 2-Monochloronaphthalene | 2 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2620 1,2-Dichloronaphthalene | 3 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2621 1,4-Dichloronaphthalene | 5 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2622 1,5-Dichloronaphthalene | 6 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2623 1,8-Dichloronaphthalene | 9 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2624 2,3-Dichloronaphthalene | 10 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2630 1,2,3-Trichloronaphthalene | 13 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2632 1,2,4-Trichloronaphthalene | 14 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2631 1,4,6-Trichloronaphthalene | 24 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2640 1,2,3,4-Tetrachloronaphthalene | 27 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2645 1,2,3,5-Tetrachloronaphthalene | 28 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2646 1,2,3,8-Tetrachloronaphthalene | 31 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2642 1,2,5,6-Tetrachloronaphthalene | 36 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2641 1,3,5,7-Tetrachloronaphthalene | 42 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2644 1,4,5,8-Tetrachloronaphthalene | 46 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2643 2,3,6,7-Tetrachloronaphthalene | 48 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2654 1,2,3,4,5-Pentachloronaphthalene | 49 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2652 1,2,3,4,6-Pentachloronaphthalene | 50 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2651 1,2,3,5,7-Pentachloronaphthalene | 52 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2650 1,2,3,5,8-Pentachloronaphthalene | 53 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2653 1,2,3,6,7-Pentachloronaphthalene | 54 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2656 1,2,4,5,8-Pentachloronaphthalene | 59 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2655 1,2,4,6,7-Pentachloronaphthalene | 60 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2668 1,2,3,4,5,6-Hexachloronaphthalene | 63 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2669 1,2,3,4,5,7-Hexachloronaphthalene | 64 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| NEW | ECN-2667 1,2,3,4,5,8-Hexachloronaphthalene | 65 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2660 1,2,3,4,6,7-Hexachloronaphthalene | 66 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2663 1,2,3,5,6,7-Hexachloronaphthalene | 67 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2664 1,2,3,5,6,8-Hexachloronaphthalene | 68 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2662 1,2,3,5,7,8-Hexachloronaphthalene | 69 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2665 1,2,3,6,7,8-Hexachloronaphthalene | 70 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2666 1,2,4,5,6,8-Hexachloronaphthalene | 71 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2661 1,2,4,5,7,8-Hexachloronaphthalene | 72 | 100 $\mu\text{g/mL}$ in nonane | Inquire |
| | ECN-2670 1,2,3,4,5,6,7-Heptachloronaphthalene | 73 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2671 1,2,3,4,5,6,8-Heptachloronaphthalene | 74 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |
| | ECN-2680 Octachloronaphthalene | 75 | 100 $\mu\text{g/mL}$ in nonane | 1 mL |

Polychlorinated Naphthalene (PCN) Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|-------------------------------------|----------------------|
| NEW ECN-5489 | PCN Calibration Solutions [CS1-CS7] | 7 x 0.5 mL in nonane |
| NEW ECN-5489-CS1 | PCN Calibration Solution [CS1] | 0.5 mL in nonane |
| NEW ECN-5489-CS2 | PCN Calibration Solution [CS2] | 0.5 mL in nonane |
| NEW ECN-5489-CS3 | PCN Calibration Solution [CS3] | 0.5 mL in nonane |
| NEW ECN-5489-CS4 | PCN Calibration Solution [CS4] | 0.5 mL in nonane |
| NEW ECN-5489-CS5 | PCN Calibration Solution [CS5] | 0.5 mL in nonane |
| NEW ECN-5489-CS6 | PCN Calibration Solution [CS6] | 0.5 mL in nonane |
| NEW ECN-5489-CS7 | PCN Calibration Solution [CS7] | 0.5 mL in nonane |

| <i>All concentrations are in ng/mL</i> | | | | | | | | |
|--|-----|-----|-----|-----|-----|-----|-----|-----|
| Unlabeled | PCN | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 |
| 1,2,3,4-TetraCN | 27 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| 1,2,3,5,7-PentaCN | 52 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| 1,2,3,4,6,7-HexaCN | 66 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| 1,2,3,5,6,7-HexaCN | 67 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| 1,2,3,5,6,8-HexaCN | 68 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| 1,2,3,4,5,6,7-HeptaCN | 73 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| OctaCN | 75 | 0.1 | 0.2 | 1 | 2 | 10 | 20 | 100 |
| Labeled | | | | | | | | |
| 1,2,3,4-TetraCN (¹³ C ₁₀ , 99%) | 27 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,5,7-PentaCN (¹³ C ₁₀ , 99%) | 52 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,5,7-HexaCN (¹³ C ₁₀ , 99%) | 64 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,5,6,7-HexaCN (¹³ C ₁₀ , 99%) | 67 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4,5,6,7-HeptaCN (¹³ C ₁₀ , 99%) | 73 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| OctaCN (¹³ C ₁₀ , 99%) | 75 | 10 | 10 | 10 | 10 | 10 | 10 | 10 |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |

| | | |
|------------------------------|-----------------------------------|----------------------------|
| NEW ECN-5490 | PCN Cleanup Solution | 5 mL in methanol/isooctane |
| NEW ECN-5490-200X-1.2 | PCN Cleanup Solution (200X stock) | Inquire |

| Labeled | PCN | ECN-5490 (ng/mL) | ECN-5490-200X-1.2 (ng/mL) |
|--|-----|---------------------|------------------------------|
| 1,2,3,4-TetraCN (¹³ C ₁₀ , 99%) | 27 | 0.5 | 100 |
| 1,2,3,5,7-PentaCN (¹³ C ₁₀ , 99%) | 52 | 0.5 | 100 |
| 1,2,3,4,5,7-HexaCN (¹³ C ₁₀ , 99%) | 64 | 0.5 | 100 |
| 1,2,3,5,6,7-HexaCN (¹³ C ₁₀ , 99%) | 67 | 0.5 | 100 |
| 1,2,3,4,5,6,7-HeptaCN (¹³ C ₁₀ , 99%) | 73 | 0.5 | 100 |
| OctaCN (¹³ C ₁₀ , 99%) | 75 | 0.5 | 100 |

| NEW ECN-5497 | PCN Native PAR Solution | 1.2 mL in nonane |
|-----------------------|-------------------------|------------------|
| Unlabeled | PCN | (ng/mL) |
| 1,2,3,4-TetraCN | 27 | 1000 |
| 1,2,3,5,7-PentaCN | 52 | 1000 |
| 1,2,3,4,6,7-HexaCN | 66 | 1000 |
| 1,2,3,5,6,7-HexaCN | 67 | 1000 |
| 1,2,3,5,6,8-HexaCN | 68 | 1000 |
| 1,2,3,4,5,6,7-HeptaCN | 73 | 1000 |
| OctaCN | 75 | 1000 |

Polychlorinated Naphthalene (PCN) Standard Mixtures

| Catalog No. | Compound | | Amount |
|-------------|--|-----|--------------------|
| ECN-5102 | Tetra-Octa PCN Mixture | | 1.2 mL in isoctane |
| | Labeled | PCN | (ng/mL) |
| | 1,2,3,4-TetraCN (¹³ C ₁₀ , 99%) | 27 | 1000 |
| | 1,3,5,7-TetraCN (¹³ C ₁₀ , 99%) | 42 | 1000 |
| | 1,2,3,5,7-PentaCN (¹³ C ₁₀ , 99%) | 52 | 1000 |
| | 1,2,3,5,6,7-HexaCN (¹³ C ₁₀ , 99%) | 67 | 1000 |
| | 1,2,3,4,5,6,7-HeptaCN (¹³ C ₁₀ , 99%) | 73 | 1000 |
| | OctaCN (¹³ C ₁₀ , 99%) | 75 | 1000 |

| | | | |
|------------|-----------------------|------------------------------|------------------|
| NEW | ECN-5558 | Mono-Octa PCN Native Mixture | 1.2 mL in nonane |
| | Unlabeled | PCN | (ng/mL) |
| | 2-MonoCN | 2 | 1000 |
| | 1,2-DiCN | 3 | 1000 |
| | 1,4-DiCN | 5 | 1000 |
| | 1,2,3-TriCN | 13 | 1000 |
| | 1,4,6-TriCN | 24 | 1000 |
| | 1,4,5,8-TetraCN | 46 | 1000 |
| | 1,3,5,7-TetraCN | 42 | 1000 |
| | 1,2,3,5,7-PentaCN | 52 | 1000 |
| | 1,2,3,5,8-PentaCN | 53 | 1000 |
| | 1,2,3,4,6,7-HexaCN | 66 | 1000 |
| | 1,2,3,5,6,8-HexaCN | 68 | 1000 |
| | 1,2,3,4,5,6,7-HeptaCN | 73 | 1000 |
| | OctaCN | 75 | 1000 |

Halowax Technical Mixtures

| Catalog No. | Compound | Concentration | Amount |
|-------------|--------------|---------------------|--------|
| ECN-1000 | HALOWAX 1000 | 100 µg/mL in hexane | 2 mL |
| ECN-1013 | HALOWAX 1013 | 100 µg/mL in hexane | 2 mL |
| ECN-1051 | HALOWAX 1051 | 100 µg/mL in hexane | 2 mL |

Priority Pollutant, Endocrine Disruptor, and Chemical Contaminant Standards

Major improvements in air and water quality resulted from focus on the prevention and remediation of priority pollutants.

Cleaning up the environment and the products that impact dietary intake will ultimately lead to a cleaner, healthier life for all of us, and future generations. Emerging contaminants, particularly those analyzed by LC-MS, are among the most active areas of CIL's new-product development efforts.



Pharmaceutical and Personal Care Product Standards

Concern about environmental and human exposure to pharmaceuticals and personal care products (PPCPs) has grown significantly. This classification encompasses a broad range of chemicals, ranging from antibiotics to hormones to pesticides. One common theme among these groups is the need for high-quality isotopically labeled standards to strengthen the analysis of PPCPs in difficult matrices such as sewage sludge and wastewater. CIL, with guidance from leading laboratories around the world, works diligently to produce representative standards for the analysis of PPCPs.

Food and Drinking Water Analysis Standards

Increased attention to possible contamination of food and water has caused analysts to broaden the scope of trace food and water testing by IDMS. Of particular interest are veterinary antibiotics used to improve the health of feed animals, ranging from shrimp to poultry to cattle. Human antibiotics, pharmaceuticals, and hormones that are not removed during wastewater treatment are also of interest, as is the routine analysis of POPs, pesticides, and other industrial contaminants that have entered the food and water supply.

Phthalate and Phthalate Metabolite Standards

Phthalates continue to be a growing environmental concern, especially as more is learned about the effect of continued exposure on the environment and the human body. Phthalate diesters are ubiquitous in the laboratory environment, so many analysts are now examining phthalate monoesters and metabolites of phthalate monoesters to reduce background interferences. Adipate esters are also anticipated to be of interest to exposure analysts; please inquire if you are interested in additional adipate standards.

Perfluorinated Compound Standards

From stain-resistant textiles to nonstick surface coatings and much more, poly- and perfluorinated compounds (PFCs) are nearly ubiquitous chemicals in the environment. CIL offers several new labeled and unlabeled perfluorinated carboxylic acid standards (PFCAs) in this catalog. CIL will be continuously adding to our offerings, so we recommend visiting our website for product updates in this rapidly growing field.

Nitrosamine Standards

Nitrosamine compounds are contaminants that may be found in food and tobacco products, and some have been classified as carcinogenic. While efforts have been made to reduce the levels of nitrosamines in commercial products, the need to monitor trace levels of this pollutant has prompted CIL to expand its offerings of labeled and unlabeled nitrosamine standards.

Halogenated and Substituted Benzene and Phenol Standards

Many industrial and consumer products are composed of chemicals that contain halogenated or substituted benzene or phenol functional groups. Resistant to decomposition and metabolism, these chemicals may persist even after the parent molecule has undergone partial decomposition, or they may exist as a product or an industrial byproduct. The increased use of brominated compounds is expected to lead to more brominated benzenes and phenols in the environment, and the continued presence of chlorinated compounds ensures that chlorinated benzenes and phenols will be found in the environment for years to come.

Bisphenol Standards

Bisphenol A (BPA) is a synthetic compound that has long been used in the production of polycarbonate plastics and epoxy resins. With recent bans on the use of BPA in certain food and water containers, replacement materials, many of which are alternative bisphenol compounds, are finding larger use. As the list of replacements grows, so too does the list of analytical standards being produced by CIL.

Perfluorokerosene Standards

Mass spectrometers require a reference compound to accurately assign masses and to verify tuning and operating conditions of the instrument. In the late 1960s, Columbia Organic Chemical Company successfully synthesized perfluorokerosene (PFK), and in a short time PFK became the most widely used reference compound in the mass spec community. Because PFK is difficult to synthesize and purify, the last producer halted production of it in 2011. CIL recognized the need for continued production of PFK and has partnered with a new producer, offering low- and high-boiling PFK standards.

Chlorinated Paraffin Standards

Chlorinated paraffins, or chloroalkanes, are industrial chemicals that have been used for many years as coolants, lubricants, plasticizers, and flame retardants. Short chain ($C_{10}-C_{13}$) chlorinated paraffins (SCCP) have come under increased scrutiny in recent years because of concerns about long-range transport, persistence in the environment, bioaccumulation, and potentially toxic endpoints. In 2006, the European Commission submitted a formal application to include SCCP in the Stockholm Convention. CIL has worked diligently to synthesize single-isomer SCCP standards to assist researchers in what is an extremely difficult analytical process.

Personal Care Product Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|-----------------------------|--------------------------------------|--------|
| DLM-183-1.2 | Benzophenone (D_{10} , 98%) | $C_6D_5COC_6D_5$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-8303-1.2 | Benzophenone (unlabeled) | $C_6H_5COCH_3$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW CLM-9437-1.2 | Decamethylcyclopentasiloxane "D5" (decamethyl- $^{13}\text{C}_{10}$, 98%) | * $C_{10}H_{30}O_5Si_5$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9442-1.2 | Decamethylcyclopentasiloxane "D5" (unlabeled) | $C_{10}H_{30}O_5Si_5$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-4762-1.2 | <i>N,N</i> -Diethyl- <i>m</i> -toluamide (DEET) | $CH_3C_6H_4CON(CH_2CD_3)_2$ | 100 $\mu\text{g/mL}$ in MeCl | 1.2 mL |
| DLM-4762-D-1.2 | (dimethyl- D_6 , 98%) | | 100 $\mu\text{g/mL}$ in dioxane | 1.2 mL |
| ULM-7975-1.2 | <i>N,N</i> -Diethyl- <i>m</i> -toluamide (DEET) (unlabeled) | $CH_3C_6H_4CON(CH_2CH_3)_2$ | 100 $\mu\text{g/mL}$ in MeCl | 1.2 mL |
| ULM-7975-D-1.2 | | | 100 $\mu\text{g/mL}$ in dioxane | 1.2 mL |
| NEW CLM-9438-1.2 | Dodecamethylcyclohexasiloxane "D6" (dodecamethyl- $^{13}\text{C}_{12}$, 98%) CP 92% | * $C_{12}H_{36}O_6Si_6$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9443-1.2 | Dodecamethylcyclohexasiloxane "D6" (unlabeled) | $C_{12}H_{36}O_6Si_6$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CLM-9349-1.2 | 4-Dodecylbenzenesulfonate, sodium salt (ring- $^{13}\text{C}_6$, 99%) CP 94% | * $C_6C_{12}H_{29}NaO_3S$ | 10 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9350-1.2 | 4-Dodecylbenzenesulfonate, sodium salt (unlabeled) | $C_{18}H_{29}NaO_3S$ | 10 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| CLM-8008-1.2 | Hexachlorophene ($^{13}\text{C}_{13}$, 99%) | * $CH_2[*C_6H(Cl)_3OH]_2$ | 50 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ULM-8009-1.2 | Hexachlorophene (unlabeled) | $CH_2[C_6H(Cl)_3OH]_2$ | 50 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CLM-9542-1.2 | Hexamethylcyclotrisiloxane "D3" (hexamethyl- $^{13}\text{C}_6$, 98%) | * $C_6H_{18}O_3Si_3$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9687-1.2 | Hexamethylcyclotrisiloxane "D3" (unlabeled) | $C_6H_{18}O_3Si_3$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| CLM-4745-1.2 | 4-Hydroxybenzoic acid (ring- $^{13}\text{C}_6$, 99%) | * $C_6CH_6O_3$ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8251-1.2 | 4-Hydroxybenzoic acid (unlabeled) | $C_6H_6O_3$ | 1 mg/mL in methanol | 1.2 mL |
| CLM-7885-1.2 | Methyl triclosan (2,4,4-trichloro-2-methoxydiphenyl ether) (ring- $^{13}\text{C}_{12}$, 99%) | * $C_{12}CH_9Cl_3O_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-7884-1.2 | Methyl triclosan (2,4,4-trichloro-2-methoxydiphenyl ether) (unlabeled) | $C_{12}CH_9Cl_3O_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW CLM-9436-1.2 | Octamethylcyclotetrasiloxane "D4" (octamethyl- $^{13}\text{C}_8$, 98%) | * $C_8H_{24}O_4Si_4$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9441-1.2 | Octamethylcyclotetrasiloxane "D4" (unlabeled) | $C_8H_{24}O_4Si_4$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CLM-9849-1.2 | Benzyl paraben (benzyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_8H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9850-1.2 | Benzyl paraben (benzyl 4-hydroxybenzoate) (unlabeled) | $C_{14}H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| CLM-8285-1.2 | <i>n</i> -Butyl paraben (ring- $^{13}\text{C}_6$, 99%) | $HO*C_6H_4CO_2(CH_2)_3CH_3$ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8287-1.2 | <i>n</i> -Butyl paraben (unlabeled) | $HOC_6H_4CO_2(CH_2)_3CH_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9761-1.2 | Ethyl paraben (ethyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_3H_{10}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9760-1.2 | Ethyl paraben (ethyl 4-hydroxybenzoate) (unlabeled) | $C_9H_{10}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9847-1.2 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_5H_{14}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9848-1.2 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (unlabeled) | $C_{11}H_{14}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9845-1.2 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_4H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9846-1.2 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (unlabeled) | $C_{10}H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| CLM-8249-1.2 | Methyl paraben (methyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_2H_8O_3$ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8250-1.2 | Methyl paraben (methyl 4-hydroxybenzoate) (unlabeled) | $C_8H_8O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9763-1.2 | <i>n</i> -Propyl paraben (<i>n</i> -propyl 4-hydroxybenzoate) (ring- $^{13}\text{C}_6$, 99%) | * $C_6C_4H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9762-1.2 | <i>n</i> -Propyl paraben (<i>n</i> -propyl 4-hydroxybenzoate) (unlabeled) | $C_{10}H_{12}O_3$ | 1 mg/mL in methanol | 1.2 mL |
| CLM-8525-1.2 | Oxybenzone (phenyl- $^{13}\text{C}_6$, 99%) | $HOC_6H_3(OCH_3)CO*C_6H_5$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-8531-1.2 | Oxybenzone (unlabeled) | $HOC_6H_3(OCH_3)COC_6H_5$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |

(continued on next page)

Personal Care Product Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------|---|--|---------------------------|--------|
| CLM-7286-1.2 | Triclocarban (3,4,4'-trichlorocarbanilide) (4'-chlorophenyl- ¹³ C ₆ , 99%) | *C ₆ C ₇ H ₉ Cl ₃ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7968-1.2 | Triclocarban (3,4,4'-trichlorocarbanilide) (unlabeled) | C ₁₃ H ₉ Cl ₃ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-6779-1.2 | Tricosan (2',4,4'-trichloro-2-hydroxydiphenyl ether) | *C ₁₂ H ₇ Cl ₃ O ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-6779-MT-1.2 | (¹³ C ₁₂ , 99%) | | 100 µg/mL in MTBE | 1.2 mL |
| ULM-6935-1.2 | Tricosan (2',4,4'-trichloro-2-hydroxydiphenyl ether) | C ₁₂ H ₇ Cl ₃ O ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-6935-MT-1.2 | (unlabeled) | | 100 µg/mL in MTBE | 1.2 mL |

Sex and Steroidal Hormone Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------|---|---|--------------------------|--------|
| NEW DLM-8438-0.001 | Aldosterone (2,2,4,6,6,17,21,21-D ₈) | C ₂₁ D ₈ H ₂₀ O ₅ | neat | 1 mg |
| NEW ULM-9134-0.001 | Aldosterone (unlabeled) CP 95% | C ₂₁ H ₂₈ O ₅ | neat | 1 mg |
| NEW ULM-9163-0.001 | 3-α,5-β-Tetrahydroaldosterone (unlabeled) | C ₂₁ H ₃₂ O ₅ | neat | 1 mg |
| NEW CLM-9135-C | 4-Androstene-3,17-dione (2,3,4- ¹³ C ₃ , 98%) | *C ₃ C ₁₆ H ₂₆ O ₂ | 100 µg/mL in methanol | 1 mL |
| NEW CLM-9135-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW DLM-8330-0.05 | 4-Androstene-3,17-dione (2,2,4,6,6-D ₅ , 98%) | C ₁₉ D ₅ H ₂₁ O ₂ | neat | 0.05 g |
| NEW ULM-8472-C | 4-Androstene-3,17-dione (unlabeled) | C ₁₉ H ₂₆ O ₂ | 100 µg/mL in methanol | 1 mL |
| NEW ULM-8472-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW DLM-9137-0.001 | Androsterone glucuronide (2,2,4,4-D ₄ , 98%) | C ₂₅ D ₄ H ₃₄ O ₈ | neat | 1 mg |
| NEW ULM-9138-0.005 | Androsterone glucuronide (unlabeled) | C ₂₅ H ₃₈ O ₈ | neat | 5 mg |
| NEW DLM-9541-0.01 | Chenodeoxycholic acid (2,2,3,4,4,6,6,7,8-D ₉ , 98%) | C ₂₄ D ₉ H ₃₁ O ₄ | neat | 10 mg |
| NEW ULM-9540-0.05 | Chenodeoxycholic acid (unlabeled) | C ₂₄ H ₄₀ O ₄ | neat | 50 mg |
| NEW DLM-8276-0.1 | Cholestenone (2,2,4,6,6-D ₅ , 98%) | C ₂₇ D ₅ H ₃₉ O | neat | 0.1 g |
| NEW CLM-9139-B | Cholesterol (2,3,4- ¹³ C ₃ , 98%) | *C ₃ C ₂₄ H ₄₆ O | 50 µg/mL in chloroform | 1 mL |
| NEW CLM-9139-C | | | 100 µg/mL in chloroform | 1 mL |
| NEW CLM-9587-1.2 | Cholesterol (23,24,25,26,27- ¹³ C ₅ , 99%) | *C ₅ C ₂₂ H ₄₆ O | 100 µg/mL in methanol | 1.2 mL |
| CLM-804-0.1 | Cholesterol (3,4- ¹³ C ₂ , 99%) | *C ₂ C ₂₅ H ₄₆ O | neat | 0.1 g |
| DLM-2607-0.1 | Cholesterol (2,2,3,4,4,6-D ₆ , 97-98%) | C ₂₇ H ₄₀ D ₆ O | neat | 0.1 g |
| DLM-3057-0.01 | Cholesterol (25,26,26,26,27,27,27-D ₇ , 98%) | C ₂₇ H ₃₉ D ₇ O | neat | 0.01 g |
| NEW ULM-9140-C | Cholesterol (unlabeled) | C ₂₇ H ₄₆ O | 100 µg/mL in chloroform | 1 mL |
| NEW ULM-9140-D | | | 1000 µg/mL in chloroform | 1 mL |
| NEW DLM-2611-0.05 | Cholic acid (2,2,4,4-D ₄ , 98%) | C ₂₄ H ₃₆ D ₄ O ₅ | neat | 50 mg |
| NEW ULM-9543-0.05 | Cholic acid (unlabeled) | C ₂₄ H ₄₀ O ₅ | neat | 50 mg |
| NEW DLM-7347-0.01 | Corticosterone (2,2,4,6,6,17α,21,21-D ₈ , 97-98%) | C ₂₁ D ₈ H ₂₂ O ₄ | neat | 0.01 g |
| NEW DLM-2057-0.01 | Cortisol (9,12,12-D ₃ , 98%) | C ₂₁ H ₂₇ D ₃ O ₅ | neat | 0.01 g |
| DLM-2218-0.1MG | Cortisol (9,11,12,12-D ₄ , 98%) | C ₂₁ D ₄ H ₂₆ O ₅ | neat | 0.1 mg |
| NEW ULM-7823-0.1MG | Cortisol (unlabeled) | C ₂₁ H ₃₀ O ₅ | neat | 0.1 mg |
| NEW ULM-9141-C | Cortisol (unlabeled) | C ₂₁ H ₃₀ O ₅ | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9141-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW DLM-9142-0.001 | Cortisone (2,2,4,6,6,12,12-D ₇ , 98%) | C ₂₁ D ₇ H ₂₁ O ₅ | neat | 1 mg |
| NEW ULM-9202-0.001 | Cortisone (unlabeled) | C ₂₁ H ₂₈ O ₅ | neat | 1 mg |
| DLM-8049-0.005 | Dehydroepiandrosterone (DHEA) (2,2,3,4,4,6-D ₆ , 99%) CP 97% | C ₁₉ H ₂₂ D ₆ O ₂ | neat | 5 mg |
| NEW ULM-9143-C | Dehydroepiandrosterone (DHEA) (unlabeled) | C ₁₉ H ₂₈ O ₂ | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9143-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW ULM-9144-C | Dehydroepiandrosterone sulfate, sodium salt (DHEAS) | C ₁₉ H ₂₈ O ₅ S | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9144-D | (unlabeled) | | 1000 µg/mL in methanol | 1 mL |
| NEW DLM-2824-0.01 | Deoxycholic acid (2,2,4,4-D ₄ , 98%) | C ₂₄ H ₃₆ D ₄ O ₄ | neat | 10 mg |
| NEW DLM-9546-0.01 | Deoxycholic acid (2,2,4,4,11,11-D ₆ , 98%) | C ₂₄ H ₃₄ D ₆ O ₄ | neat | 10 mg |
| NEW ULM-9545-0.05 | Deoxycholic acid (unlabeled) | C ₂₄ H ₄₀ O ₄ | neat | 50 mg |
| NEW DLM-8305-0.01 | 21-Deoxycortisol (2,2,4,6,6,21,21,21-D ₈ , 97%) | C ₂₁ D ₈ H ₂₂ O ₄ | neat | 0.01 g |
| NEW ULM-9145-C | 11-Deoxycortisol (unlabeled) | C ₂₁ H ₃₀ O ₄ | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9145-D | | | 1000 µg/mL in methanol | 1 mL |
| DLM-170-D-1.2 | Diethylstilbestrol (<i>cis/trans</i> mix) (ring-3,3',5,5'-diethyl-1,1',1'-D ₈ , 98%) | HOC ₆ D ₂ H(CH ₂ CD ₂) C=C(CD ₂ CH ₃)C ₆ H ₂ D ₂ OH | 100 µg/mL in dioxane | 1.2 mL |

Sex and Steroidal Hormone Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------------|---|---------------------------------|---------------------------|--------|
| ULM-7921-D-1.2 | Diethylstilbestrol (<i>cis/trans</i> mix) (unlabeled) | <chem>OCC6C(C)C(O)C(C)C6</chem> | 100 µg/mL in dioxane | 1.2 mL |
| NEW CLM-9146-C | 5-α-Dihydrotestosterone (2,3,4- ¹³ C ₃ , 99%) | * <chem>C3C16H28O2</chem> | 100 µg/mL in methanol | 1 mL |
| NEW CLM-9146-D | | | 1000 µg/mL in methanol | 1 mL |
| CLM-7936-0.1MG | DL-Estradiol (13,14,15,16,17,18- ¹³ C ₆ , 99%) | <chem>C12C6H24O2</chem> | neat | 0.1 mg |
| CLM-7936-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| CLM-803-1.2 | Estradiol (3,4- ¹³ C ₂ , 99%) | * <chem>C2C16H24O2</chem> | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-2487-5 | Estradiol (2,4,16,16-D ₄ , 95-97%) | <chem>C18H20D4O2</chem> | neat | 5 mg |
| NEW ULM-7449-0.1MG | Estradiol (unlabeled) | <chem>C18H24O2</chem> | neat | 0.1 mg |
| ULM-7449-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-9147-C | Estriol (16-α-hydroxyestradiol) (2,3,4- ¹³ C ₃ , 99%) | * <chem>C3C15H24O3</chem> | 100 µg/mL in methanol | 1 mL |
| DLM-8583-0.1MG | Estriol (2,4,16,17-D ₄ , 98%) CP 95% | <chem>C18D4H20O3</chem> | neat | 0.1 mg |
| ULM-8218-0.1MG | Estriol (unlabeled) | <chem>C18H24O3</chem> | neat | 0.1 mg |
| CLM-7935-0.1MG | DL-Estrone (13,14,15,16,17,18- ¹³ C ₆ , 99%) | <chem>C12C6H22O2</chem> | neat | 0.1 mg |
| CLM-7935-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-9148-B | Estrone (2,3,4- ¹³ C ₃ , 99%) | * <chem>C3C15H22O2</chem> | 50 µg/mL in methanol | 1 mL |
| NEW CLM-9148-C | | | 100 µg/mL in methanol | 1 mL |
| CLM-673-1.2 | Estrone (3,4- ¹³ C ₂ , 90%) | * <chem>C2C16H22O2</chem> | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-3976-5 | Estrone (2,4,16,16-D ₄ , 97%) | <chem>C18H18D4O2</chem> | neat | 5 mg |
| ULM-7212-1.2 | Estrone (unlabeled) | <chem>C18H22O2</chem> | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-8033-0.1MG | DL-Estrone 3-methyl ether (13,14,15,16,17,18- ¹³ C ₆ , 99%) | * <chem>C6C13H24O2</chem> | neat | 0.1 mg |
| NEW CLM-8018-0.1MG | DL-Sodium estrone 3-sulfate (13,14,15,16,17,18- ¹³ C ₆ , 99%) (12-13%-sodium acetate) | * <chem>C6C12H21O5SNa</chem> | neat | 0.1 mg |
| NEW ULM-8132-0.1MG | Sodium estrone 3-sulfate (unlabeled) | <chem>C18H21O5SNa</chem> | neat | 0.1 mg |
| CLM-3375-1.2 | Ethyneestradiol (20,21- ¹³ C ₂ , 99%) | * <chem>C2C18H24O2</chem> | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-4691-0.01 | 17-α-Ethyneestradiol (2,4,16,16-D ₄ , 97-98%) | <chem>C20D4H20O2</chem> | neat | 0.01 g |
| ULM-7211-1.2 | Ethyneestradiol (unlabeled) | <chem>C20H24O2</chem> | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-9550-0.01 | Glycochenodeoxycholic acid (2,2,3,4,4,6,6,7,8-D ₉ , 98%) CP 97% | <chem>C26D9H34NO5</chem> | neat | 10 mg |
| NEW DLM-2742-0.01 | Glycocholic acid (2,2,4,4-D ₄ , 98%) (contains ~4% water) | <chem>C26D4H39NO6</chem> | neat | 10 mg |
| NEW ULM-9551-0.05 | Glycocholic acid (unlabeled) | <chem>C26H43NO6</chem> | neat | 50 mg |
| NEW DLM-9553-0.01 | Glycodeoxycholic acid (2,2,4,4,11,11-D ₆ , 98%) | <chem>C26D6H37NO5</chem> | neat | 10 mg |
| NEW DLM-9554-0.01 | Glycodeoxycholic acid (2,2,4,4-D ₄ , 98%) | <chem>C26D4H39NO5</chem> | neat | 10 mg |
| NEW ULM-9552-0.05 | Glycodeoxycholic acid, sodium salt (unlabeled) | <chem>C26H42NNaO5</chem> | neat | 50 mg |
| NEW DLM-9556-0.01 | Glycolithocholic acid (2,2,4,4-D ₄ , 98%) | <chem>C26D4H39NO4</chem> | neat | 10 mg |
| NEW ULM-9555-0.05 | Glycolithocholic acid, sodium salt (unlabeled) | <chem>C26H42NNaO4</chem> | neat | 50 mg |
| NEW DLM-9558-0.01 | Glycoursoodeoxycholic acid (2,2,4,4-D ₄ , 98%) CP 97% | <chem>C26D4H39NO5</chem> | neat | 10 mg |
| NEW ULM-9557-0.05 | Glycoursoodeoxycholic acid (unlabeled) | <chem>C26H43NO5</chem> | neat | 50 mg |
| NEW DLM-9150-0.001 | 18-Hydroxycorticosterone (9,11,12,12-D ₄ , 98%) CP 95% | <chem>C21D4H26O5</chem> | neat | 1 mg |
| NEW ULM-9151-0.001 | 18-Hydroxycorticosterone (unlabeled) CP 95% | <chem>C21H30O5</chem> | neat | 1 mg |
| NEW DLM-9149-0.001 | 6-β-Hydroxycortisol (9,11,12,12-D ₄ , 98%) CP 97% | <chem>C21D4H26O6</chem> | neat | 1 mg |
| CLM-8012-0.1MG | DL-2-Hydroxyestradiol (13,14,15,16,17,18- ¹³ C ₆ , 99%) | * <chem>C6C12H24O3</chem> | neat | 0.1 mg |
| ULM-8135-0.1MG | 2-Hydroxyestradiol (unlabeled) | <chem>C18H24O3</chem> | neat | 0.1 mg |
| NEW CLM-9153-0.1MG | 16-α-Hydroxyestrone (2,3,4- ¹³ C ₃ , 99%) | * <chem>C3C15H22O3</chem> | neat | 0.1 mg |
| CLM-8011-0.1MG | DL-2-Hydroxyestrone (13,14,15,16,17,18- ¹³ C ₆ , 99%) | * <chem>C6C12H22O3</chem> | neat | 0.1 mg |
| ULM-8134-0.1MG | 2-Hydroxyestrone (unlabeled) | <chem>C18H22O3</chem> | neat | 0.1 mg |
| CLM-8013-0.1MG | DL-4-Hydroxyestrone (13,14,15,16,17,18- ¹³ C ₆ , 99%) | * <chem>C6C12H22O3</chem> | neat | 0.1 mg |
| ULM-8261-0.1MG | 4-Hydroxyestrone (unlabeled) CP 96% | <chem>C18H22O3</chem> | neat | 0.1 mg |
| CLM-8016-0.1MG | DL-2-Hydroxyestrone-3-methyl ether (13,14,15,16,17,18- ¹³ C ₆ , 99%) | <chem>C13C6H24O3</chem> | neat | 0.1 mg |
| ULM-8133-0.1MG | 2-Hydroxyestrone-3-methyl ether (unlabeled) | <chem>C19H24O3</chem> | neat | 0.1 mg |
| NEW CDLM-9154-C | 17α-Hydroxypregnanolone | * <chem>C2C19D2H30O3</chem> | 100 µg/mL in methanol | 1 mL |
| NEW CDLM-9154-D | (20,21- ¹³ C ₂ , 99%; 16,16-D ₂ , 99%) | | 1000 µg/mL in methanol | 1 mL |
| NEW ULM-9155-C | 17α-Hydroxypregnanolone (unlabeled) | <chem>C21H32O3</chem> | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9155-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW CLM-9157-C | 17α-Hydroxyprogesterone (2,3,4- ¹³ C ₃ , 98%) | * <chem>C3C18H30O3</chem> | 100 µg/mL in methanol | 1 mL |
| NEW CLM-9157-D | | | 1000 µg/mL in methanol | 1 mL |

(continued on next page)

Sex and Steroidal Hormone Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------|---|---|---------------------------|--------|
| NEW DLM-6598-0.01 | 17-Hydroxyprogesterone (2,2,4,6,6,21,21,21-D ₈ , 98%) | C ₂₁ H ₂₂ D ₈ O ₃ | neat | 0.01 g |
| NEW ULM-9156-C | 17α-Hydroxyprogesterone (unlabeled) CP 95% | C ₂₁ H ₃₀ O ₃ | 100 µg/mL in methanol | 1 mL |
| NEW ULM-9156-D | | | 1000 µg/mL in methanol | 1 mL |
| NEW DLM-9560-0.05 | Lithocholic acid (2,2,4,4-D ₄ , 98%) | C ₂₄ D ₄ H ₃₆ O ₃ | neat | 50 mg |
| NEW ULM-9559-0.05 | Lithocholic acid (unlabeled) | C ₂₄ H ₄₀ O ₃ | neat | 50 mg |
| CLM-8015-0.1MG | DL-2-Methoxyestradiol (13,14,15,16,17,18- ¹³ C ₆ , 99%) | *C ₆ C ₁₃ H ₂₆ O ₃ | neat | 0.1 mg |
| ULM-8137-0.1MG | 2-Methoxyestradiol (unlabeled) | C ₁₉ H ₂₆ O ₃ | neat | 0.1 mg |
| CLM-8019-0.1MG | DL-4-Methoxyestradiol (13,14,15,16,17,18- ¹³ C ₆ , 99%) | C ₁₃ *C ₆ H ₂₆ O ₃ | neat | 0.1 mg |
| NEW ULM-8136-0.1MG | 4-Methoxyestradiol (unlabeled) | C ₁₉ H ₂₆ O ₃ | neat | 0.1 mg |
| CLM-8014-0.1MG | DL-2-Methoxyestrone (13,14,15,16,17,18- ¹³ C ₆ , 99%) | *C ₆ C ₁₃ H ₂₄ O ₃ | neat | 0.1 mg |
| ULM-8263-0.1MG | 2-Methoxyestrone (unlabeled) | C ₁₉ H ₂₄ O ₃ | neat | 0.1 mg |
| CLM-8017-0.1MG | DL-4-Methoxyestrone (13,14,15,16,17,18- ¹³ C ₆ , 99%) | C ₁₃ *C ₆ H ₂₄ O ₃ | neat | 0.1 mg |
| ULM-8262-0.1MG | 4-Methoxyestrone (unlabeled) | C ₁₉ H ₂₄ O ₃ | neat | 0.1 mg |
| NEW CLM-2468-0.01 | Norethindrone (ethynodiol- ¹³ C ₂ , 99%) | *C ₂ C ₁₈ H ₂₆ O ₂ | neat | 0.01 g |
| NEW DLM-3979-1.2 | 19-Nortestosterone (16,16,17-D ₃ , 98%) | C ₁₈ H ₂₃ D ₃ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| DLM-3979-5 | | | neat | 5 mg |
| NEW ULM-4841-1.2 | 19-Nortestosterone (unlabeled) | C ₁₈ H ₂₆ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW DLM-3754-0.01 | 5-α-Pregnan-3-α-ol-20-one (17,21,21,21-D ₄ , 96-98%) | C ₂₁ H ₃₀ D ₄ O ₂ | neat | 0.01 g |
| | CP 95%+ | | | |
| NEW DLM-2294-0.01 | 5-β-Pregnan-3-α-ol-20-one (17,21,21,21-D ₄ , 96-98%) | C ₂₁ H ₃₀ D ₄ O ₂ | neat | 0.01 g |
| NEW DLM-3816-0.01 | 5-α-Pregnane-3,20-dione (1,2,4,5,6,7-D ₆ , 95%) | C ₂₁ H ₂₆ D ₆ O ₂ | neat | 0.01 g |
| NEW DLM-3910-0.01 | 5-α-Pregnane-3-α,21-diol-20-one (17,21,21-D ₃ , 95%) | C ₂₁ H ₃₁ D ₃ O ₃ | neat | 0.01 g |
| CDLM-9158-0.001 | Pregnenolone (20,21- ¹³ C ₂ , 99%; 16,16-D ₂ , 98%) | *C ₂ C ₁₉ D ₂ H ₃₀ O ₂ | neat | 1 mg |
| NEW ULM-9159-0.001 | Pregnenolone (unlabeled) | C ₂₁ H ₃₂ O ₂ | neat | 1 mg |
| NEW CDLM-9160-0.001 | Pregnenolone sulfate, sodium salt (20,21- ¹³ C ₂ , 99%; 16,16-D ₂ , 98%) | *C ₂ C ₁₉ D ₂ H ₂₉ NaO ₅ S | neat | 1 mg |
| NEW ULM-9161-0.001 | Pregnenolone sulfate, sodium salt (unlabeled) | C ₂₁ H ₃₁ NaO ₅ S | neat | 1 mg |
| NEW CLM-9162-B | Progesterone (2,3,4- ¹³ C ₃ , 99%) | *C ₃ C ₁₈ H ₃₀ O ₂ | 50 µg/mL in acetonitrile | 1 mL |
| NEW CLM-9162-C | | | 100 µg/mL in acetonitrile | 1 mL |
| NEW CLM-457-0.01 | Progesterone (3,4- ¹³ C ₂ , 90%) | *C ₂ C ₁₉ H ₃₀ O ₂ | neat | 0.01 g |
| NEW DLM-7953-1.2 | Progesterone (2,2,4,6,6,17α,21,21,21-D ₉ , 98%) | C ₂₁ D ₉ H ₂₁ O ₂ | 100 µg/mL in dioxane | 1.2 mL |
| ULM-8219-1.2 | Progesterone (unlabeled) | C ₂₁ H ₃₀ O ₂ | 100 µg/mL in dioxane | 1.2 mL |
| NEW DLM-9562-0.01 | Taurochenodeoxycholic acid, sodium salt (2,2,4,4-D ₄ , 98%) CP 97% | C ₂₆ D ₄ H ₄₀ NNaO ₆ S | neat | 10 mg |
| NEW DLM-9563-0.005 | Taurochenodeoxycholic acid, sodium salt (2,2,3,4,4,6,6,7,8-D ₉ , 98%) | C ₂₆ D ₉ H ₃₅ NNaO ₆ S | neat | 5 mg |
| NEW ULM-9561-0.05 | Taurochenodeoxycholic acid, sodium salt (unlabeled) | C ₂₆ H ₄₄ NNaO ₆ S | neat | 50 mg |
| NEW DLM-9568-0.01 | Taurodeoxycholic acid, sodium salt (2,2,4,4-D ₄ , 98%) | C ₂₆ D ₄ H ₄₀ NNaO ₆ S | neat | 10 mg |
| NEW DLM-9567-0.005 | Taurodeoxycholic acid, sodium salt (2,2,4,4,11,11-D ₆ , 98%) | C ₂₆ D ₆ H ₃₈ NNaO ₆ S | neat | 5 mg |
| NEW DLM-9570-0.01 | Taurolithocholic acid, sodium salt (2,2,4,4-D ₄ , 98%) | C ₂₆ D ₄ H ₄₀ NO ₅ SNa | neat | 10 mg |
| NEW ULM-9569-0.05 | Taurolithocholic acid, sodium salt (unlabeled) | C ₂₆ H ₄₄ NO ₅ SNa | neat | 50 mg |
| NEW CLM-159-0.01 | Testosterone (3,4- ¹³ C ₂ , 99%) | *C ₂ C ₁₇ H ₂₈ O ₂ | neat | 0.01 g |
| NEW CLM-9164-C | Testosterone (2,3,4- ¹³ C ₃ , 99%) | *C ₃ C ₁₆ H ₂₈ O ₂ | 100 µg/mL in methanol | 1 mL |
| DLM-683-1.2 | Testosterone (1,2-D ₂ , 98%) | C ₁₉ D ₂ H ₂₆ O ₂ | 100 µg/mL in MeCl | 1.2 mL |
| DLM-8085-1.2 | Testosterone (2,2,4,6,6-D ₅ , 98%) | C ₁₉ D ₅ H ₂₃ O ₂ | 100 µg/mL in MeCl | 1.2 mL |
| DLM-8085-D-1.2 | | | 100 µg/mL in dioxane | 1.2 mL |
| NEW COLM-9061-1.2 | Testosterone (3,4- ¹³ C ₂ , 99%; 17- ¹⁸ O, 98%) | *C ₂ C ₁₇ H ₂₈ *OO | 100 µg/mL in MeCl | 1.2 mL |
| ULM-8081-1.2 | Testosterone (unlabeled) | C ₁₉ H ₂₈ O ₂ | 100 µg/mL in MeCl | 1.2 mL |
| ULM-8081-D-1.2 | | | 100 µg/mL in dioxane | 1.2 mL |
| CLM-6725-0.1MG | L-Thyroxine (tyrosine-ring- ¹³ C ₆ , 99%) CP 90% | *C ₆ C ₉ H ₁₁ I ₄ NO ₄ | neat | 0.1 mg |
| NEW CLM-8931-0.1MG | L-Thyroxine (ring- ¹³ C ₁₂ , 99%) CP 97% | HO*C ₆ H ₂ (I) ₂ O*C ₆ H ₂ (I) ₂ CH ₂ CH(NH ₂)CO ₂ H | neat | 0.1 mg |
| NEW ULM-8184-0.2MG | L-Thyroxine (unlabeled) | C ₁₅ H ₁₁ I ₄ NO ₄ | neat | 0.2 mg |
| NEW DLM-9574-0.05 | Ursodeoxycholic acid (2,2,4,4-D ₄ , 98%) | C ₂₄ D ₄ H ₃₆ O ₄ | neat | 50 mg |
| NEW ULM-9573-0.05 | Ursodeoxycholic acid (unlabeled) | C ₂₄ H ₄₀ O ₄ | neat | 50 mg |

Prescription and Nonprescription Drug Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------|---|---|--|---------|
| CNLM-3726-1.2 | Acetaminophen (acetyl- ¹³ C ₂ , 99%; ¹⁵ N, 98%) | *CH ₃ *CO*NHC ₆ H ₄ OH | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7629-1.2 | Acetaminophen (unlabeled) | CH ₃ CONHC ₆ H ₄ OH | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-3008-1.2 | Amitriptyline·HCl (N,N-dimethyl-D ₆ , 98%) | C ₂₀ H ₁₇ D ₆ N·HCl | 100 µg/mL in methanol | 1.2 mL |
| ULM-8350-1.2 | Amitriptyline·HCl (unlabeled) | C ₂₀ H ₂₃ N·HCl | 100 µg/mL in methanol | 1.2 mL |
| CLM-514-1.2 | Caffeine (trimethyl- ¹³ C ₃ , 99%) | *C ₃ C ₅ H ₁₀ N ₄ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-7653-1.2 | Caffeine (unlabeled) | C ₈ H ₁₀ N ₄ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| DLM-2806-1.2 | Carbamazepine (D ₁₀ , 98%) | C ₁₅ D ₁₀ H ₂ N ₂ O | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| ULM-6581-1.2 | Carbamazepine (unlabeled) CP 97% | C ₁₅ H ₁₂ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-1287-1.2 | Clonidine (4,4,5,5-imidazoline-D ₄ , 98%) | C ₉ H ₅ D ₄ N ₃ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-8349-1.2 | Clonidine (unlabeled) | C ₉ H ₉ N ₃ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| C-041 | Codeine (D ₆ , 98%) | C ₁₈ H ₁₅ D ₆ NO ₃ | 1.0 mg/mL in methanol | 1 mL |
| C-006 | Codeine (unlabeled) | C ₁₈ H ₂₁ NO ₃ | 1.0 mg/mL in methanol | 1 mL |
| NEW DLM-1819-1.2 | DL-Cotinine (methyl-D ₃ , 98%) | C ₁₀ H ₉ D ₃ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9614-1.2 | Cotinine (unlabeled) | C ₁₀ H ₁₂ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9614-W-1.2 | | | 100 µg/mL in water | 1.2 mL |
| D-902 | Diazepam (D ₅ , 98%) | C ₁₆ H ₇ D ₅ N ₂ O·HCl | 100 µg/mL in methanol | 1 mL |
| D-907 | Diazepam (unlabeled) | C ₁₆ H ₁₂ N ₂ O·HCl | 1.0 mg/mL in methanol | 1 mL |
| NEW DLM-9974-1.2 | Diclofenac sodium (D ₄ , 98%) | C ₁₄ D ₄ H ₆ Cl ₂ NNaO ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9975-1.2 | Diclofenac sodium (unlabeled) | C ₁₄ H ₁₀ Cl ₂ NNaO ₂ | 100 µg/mL in methanol | 1.2 mL |
| CNLM-411-1.2 | 5,5-Diphenylhydantoin (2- ¹³ C, 99%; 1,3- ¹⁵ N ₂ , 98%) | *CC ₁₄ H ₁₂ *N ₂ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-8533-1.2 | 5,5-Diphenylhydantoin (unlabeled) | C ₁₅ H ₁₂ N ₂ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| F-919 | Fluoxetine oxalate (D ₆ , 98%) | C ₁₇ H ₁₂ D ₆ F ₃ NO·C ₂ H ₂ O ₄ | 100 µg/mL in methanol | 1 mL |
| F-918 | Fluoxetine-HCl (unlabeled) | C ₁₇ H ₁₈ F ₃ NO·HCl | 1.0 mg/mL in methanol | 1 mL |
| DLM-8221-1.2 | Gemfibrozil (2,2-dimethyl-D ₆ , 98%) | C ₁₅ D ₆ H ₁₆ O ₃ | 100 µg/mL in p-dioxane | 1.2 mL |
| ULM-8225-1.2 | Gemfibrozil (unlabeled) | C ₁₅ H ₂₂ O ₃ | 100 µg/mL in p-dioxane | 1.2 mL |
| CLM-6943-1.2 | Ibuprofen (propionic- ¹³ C ₃ , 99%) | *C ₃ C ₁₀ H ₁₈ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7275-1.2 | Ibuprofen (unlabeled) | C ₁₃ H ₁₈ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-3035-1.2 | Imipramine·HCl (2,4,6,8-D ₄ , 98%) | C ₁₉ H ₂₀ D ₄ N ₂ ·HCl | 100 µg/mL in methanol | 1.2 mL |
| I-902 | Imipramine (unlabeled) | C ₁₉ H ₂₄ N ₂ | 1.0 mg/mL in methanol | 1 mL |
| L-902 | Lorazepam (D ₄ , 98%) | C ₁₅ H ₆ D ₄ N ₂ O ₂ Cl ₂ | 100 µg/mL in acetonitrile | 1 mL |
| L-901 | Lorazepam (unlabeled) | C ₁₅ H ₁₀ N ₂ O ₂ Cl ₂ | 1.0 mg/mL in acetonitrile | 1 mL |
| CDLM-7665-1.2 | Naproxen (methyl- ¹³ C, 99% methyl-D ₃ , 98%) | *CC ₁₃ D ₃ H ₁₁ O ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7709-1.2 | Naproxen (unlabeled) | C ₁₄ H ₁₄ O ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CNLM-8223-1.2 | Nitrofurazone (carbonyl- ¹³ C, 99%; hydrazine- ¹⁵ N ₂ , 98%) CP 97%+ | *CC ₅ H ₆ *N ₂ N ₂ O ₄ | | Inquire |
| NEW ULM-8234 | Nitrofurazone (unlabeled) | C ₆ H ₆ N ₄ O ₄ | | Inquire |
| N-922 | Norfluoxetine oxalate (D ₆ , 98%) | C ₁₆ H ₁₀ D ₆ F ₃ NO·C ₂ H ₂ O ₄ | 100 µg/mL in methanol | 1 mL |
| N-923 | Norfluoxetine oxalate (unlabeled) | C ₁₆ H ₁₆ F ₃ NO·C ₂ H ₂ O ₄ | 1.0 mg/mL in methanol | 1 mL |
| DLM-3039-1MG | Phenylbutazone (diphenyl-D ₁₀ , 98%) | C ₁₉ D ₁₀ H ₁₀ N ₂ O ₂ | neat | 1 mg |
| NEW DLM-3039-0.05 | | | neat | 0.05 g |
| NEW DLM-3039-0.1 | | | neat | 0.1 g |
| ULM-7378-1MG | Phenylbutazone (unlabeled) | C ₁₉ H ₂₀ N ₂ O ₂ | neat | 1 mg |
| CLM-7892 | Resorcinol (¹³ C ₆ , 99%) | *C ₆ H ₆ O ₂ | | Inquire |
| CLM-8370-1.2 | Thiabendazole (ring- ¹³ C ₆ , 99%) | C ₄ *C ₆ H ₇ N ₃ S | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8371-1.2 | Thiabendazole (unlabeled) | C ₁₀ H ₇ N ₃ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-6861-MT-1.2 | Warfarin (phenyl-D ₅ , 98%) | C ₁₉ H ₁₁ D ₅ O ₄ | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-7242-MT-1.2 | Warfarin (unlabeled) | C ₁₉ H ₁₆ O ₄ | 100 µg/mL in MTBE | 1.2 mL |

Veterinary and Human Antibiotic Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|----------------------------|---|--|---------------------------|--------|
| CLM-7407-1MG | Amoxicillin-3H ₂ O (phenyl- ¹³ C ₆ , 99%) | *C ₆ C ₁₀ H ₁₉ N ₃ O ₅ S·3H ₂ O | neat | 1 mg |
| DLM-119-1.2 | (+/-)-Chloramphenicol (ring-D ₄ , benzyl-D ₁ , 98%) | NO ₂ (C ₆ D ₄)CH(OH)CH(NHCOCHCl ₂)CH ₂ OH | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-6687-1.2 | (+/-)-Chloramphenicol (unlabeled) | NO ₂ (C ₆ H ₄)CH(OH)CH(NHCOCHCl ₂)CH ₂ OH | 100 µg/mL in acetonitrile | 1.2 mL |
| CNLM-7539-1.2 | Ciprofloxacin-HCl (2,3,carboxyl- ¹³ C ₃ , 99%; quinoline- ¹⁵ N, 98%) | *C ₃ C ₁₄ H ₁₈ F*NN ₂ O ₃ ·HCl | 100 µg/mL in methanol | 1.2 mL |
| ULM-7710-1.2 | Ciprofloxacin-HCl (unlabeled) | C ₁₇ H ₁₈ FN ₃ O ₃ ·HCl | 100 µg/mL in methanol | 1.2 mL |
| <i>NEW</i> CLM-3672-MT-1.2 | Erythromycin (90-95% Erythromycin A) (N,N-dimethyl- ¹³ C ₂ , ~90%) | *C ₂ C ₃₅ H ₆₇ NO ₁₃ | 100 µg/mL in MTBE | 1.2 mL |
| <i>NEW</i> ULM-4322-MT-1.2 | Erythromycin (unlabeled) | C ₃₇ H ₆₇ NO ₁₃ | 100 µg/mL in MTBE | 1.2 mL |
| CLM-3045-1.2 | Sulfamethazine (phenyl- ¹³ C ₆ , 90%) | H ₂ N*C ₆ H ₄ SO ₂ NH(C ₆ N ₂ H ₇) | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7220-1.2 | Sulfamethazine (unlabeled) | H ₂ NC ₆ H ₄ SO ₂ NH(C ₆ N ₂ H ₇) | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-6944-1.2 | Sulfamethoxazole (ring- ¹³ C ₆ , 99%) | C ₄ *C ₆ H ₁₁ N ₃ O ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7527-1.2 | Sulfamethoxazole (unlabeled) | C ₁₀ H ₁₁ N ₃ O ₃ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7988-A-1.2 | Trimethoprim (¹³ C ₃ , 99%) | *C ₃ C ₁₁ H ₁₈ N ₄ O ₃ | 50 µg/mL in methanol | 1.2 mL |
| ULM-7989-A-1.2 | Trimethoprim (unlabeled) | C ₁₄ H ₁₈ N ₄ O ₃ | 50 µg/mL in methanol | 1.2 mL |

Tobacco-Specific Nitrosamines and Other Tobacco-Related Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-----------------------------|---|--|-----------------------------------|--------|
| <i>NEW</i> CLM-6651-1.2 | Anabasine (2,2'3,4,5,6- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₄ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-7281-1.2 | Anabasine (unlabeled) | C ₁₀ H ₁₄ N ₂ | 0.1 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-6652-1.2 | Anatabine (2,2',3,4,5,6- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₂ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-7282-1.2 | Anatabine (unlabeled) | C ₁₀ H ₁₂ N ₂ | 0.1 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-9692-1.2 | DL-Cotinine (2',3',4'- ¹³ C ₃ , 99%) CP 97% | *C ₃ C ₇ H ₁₂ N ₂ O | 100 µg/mL in water | 1.2 mL |
| <i>NEW</i> DLM-1819-1.2 | DL-Cotinine (methyl-D ₃ , 98%) | C ₁₀ D ₃ H ₉ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9614-1.2 | Cotinine (unlabeled) | C ₁₀ H ₁₂ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9614-W-1.2 | | | 100 µg/mL in water | 1.2 mL |
| CLM-6023-1.2 | 4-Methylumbelliferon (2,3,4,methyl- ¹³ C ₄ , 99%) | *C ₄ C ₆ H ₈ O ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7309-1.2 | 4-Methylumbelliferon (unlabeled) | C ₁₀ H ₈ O ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-6705-1.2 | NAB (N'-Nitrosoanabasine) (¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₃ N ₃ O | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-7168-1.2 | NAB (N'-Nitrosoanabasine) (unlabeled) | C ₁₀ H ₁₃ N ₃ O | 0.5 mg/mL in acetonitrile | 1.2 mL |
| | | | 2 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-6704-1.2 | NAT (N'-Nitrosoanatabine) (¹³ C ₆ , 99%) CP 95% | *C ₆ C ₄ H ₁₁ N ₃ O | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7207-1.2 | NAT (N'-Nitrosoanatabine) (unlabeled) | C ₁₀ H ₁₁ N ₃ O | 2 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-3914-1.2 | DL-Nicotine (3',4',5'- ¹³ C ₃ , 99%) | *C ₃ C ₇ H ₁₄ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9547-1.2 | Nicotine (unlabeled) | C ₁₀ H ₁₄ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-4556-1.2 | NNAL (4-(methylnitrosamino)-1-(3-pyridyl)-1-butanol) (1,2',3',4',5',6'- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₅ N ₃ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9434-1.2 | NNAL (4-(methylnitrosamino)-1-(3-pyridyl)-1-butanol) | C ₁₀ H ₁₅ N ₃ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9434-20X-1.2 | (unlabeled) | | 2 mg/mL in acetonitrile | 1.2 mL |
| CLM-4555-1.2 | NNK (Nicotine-derived nitrosamine ketone) | *C ₆ C ₄ H ₁₃ N ₃ O ₂ | 100 µg/mL in nonane/ethanol (9:1) | 1.2 mL |
| | (1,2',3',4',5',6'- ¹³ C ₆ , 99%) | | | |
| <i>NEW</i> ULM-8987-1.2 | NNK (Nicotine-derived nitrosamine ketone) (unlabeled) | C ₁₀ H ₁₃ N ₃ O ₂ | 100 µg/mL in nonane/ethanol (9:1) | 1.2 mL |
| <i>NEW</i> ULM-8987-20X-1.2 | | | 2 mg/mL in acetonitrile | 1.2 mL |
| CLM-4557-1.2 | NNN (N-Nitrosonornicotine) (2,2',3,4,5,6- ¹³ C ₆ , 99%) | *C ₆ C ₃ H ₁₁ N ₃ O | 100 µg/mL in nonane/ethanol (9:1) | 1.2 mL |
| <i>NEW</i> DLM-7474-1.2 | NNN (N-Nitrosonornicotine) (2,4,5,6-D ₄ , 98%) | C ₉ D ₄ H ₉ N ₃ O | 0.1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9406-1.2 | NNN (N-Nitrosonornicotine) (unlabeled) | C ₉ H ₁₁ N ₃ O | 0.1 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9406-20X-1.2 | | | 2 mg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-4896-1.2 | DL-Norcotinine (3',4',5'- ¹³ C ₃ , 99%) | *C ₃ C ₆ H ₁₀ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-9615-1.2 | Norcotinine (unlabeled) | C ₉ H ₁₀ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> CLM-4892-1.2 | DL-Nornicotine (3',4',5'- ¹³ C ₃ , 99%) | *C ₃ C ₆ H ₁₂ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| <i>NEW</i> ULM-2154-1.2 | Nornicotine (unlabeled) | C ₉ H ₁₂ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |

Food and Drinking Water Impurity Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|----------------------|--|--|---|--------|
| CLM-813-1.2 | Acrylamide (+100 ppm hydroquinone) (1,2,3- ¹³ C ₃ , 99%) | H ₂ *C=CH*CONH ₂ | 1 mg/mL in methanol | 1.2 mL |
| ULM-6721-1.2 | Acrylamide (+100 ppm hydroquinone) (unlabeled) | H ₂ C=CHCONH ₂ | 1 mg/mL in methanol | 1.2 mL |
| DLM-7170-1.2 | 1-Aminohydantoin hydrochloride (AHD) (5,5-D ₂ , 98%) | C ₃ H ₂ D ₂ N ₂ O ₂ Cl | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| ULM-7188-1.2 | 1-Aminohydantoin hydrochloride (AHD) (unlabeled) | C ₃ H ₂ N ₂ O ₂ ·HCl | 100 µg/mL in methanol | 1.2 mL |
| DLM-7171-1.2 | 3-Amino-2-oxazolidone (AOZ) (ring-D ₄ , 98%) | C ₃ H ₂ D ₄ N ₂ O ₂ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| ULM-7189-1.2 | 3-Amino-2-oxazolidone (AOZ) (unlabeled) | C ₃ H ₂ N ₂ O ₂ ·HCl | 100 µg/mL in methanol | 1.2 mL |
| DLM-7172-1.2 | 5-(4-Morpholinylmethyl)-3-amino-2-oxazolidinone (AMOZ) (4,4,5,5'-D ₅ , 98%) | C ₈ H ₁₀ D ₅ N ₃ O ₃ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| ULM-7190-1.2 | 5-(4-Morpholinylmethyl)-3-amino-2-oxazolidinone (AMOZ) (unlabeled) | C ₈ H ₁₅ N ₃ O ₃ | 100 µg/mL in methanol | 1.2 mL |
| CLM-8589-1.2 | Ammelide (ring- ¹³ C ₃ , 99%) | *C ₃ H ₄ N ₄ O ₂ | 100 µg/mL in water/diethylamine (80/20 v/v) | 1.2 mL |
| ULM-8590-1.2 | Ammelide (unlabeled) | C ₃ H ₄ N ₄ O ₂ | 100 µg/mL in water/diethylamine (80/20 v/v) | 1.2 mL |
| CLM-8316-1.2 | Ammeline (desethyldesisopropylhydroxyatrazine) (ring- ¹³ C ₃ , 99%) | *C ₃ H ₅ N ₅ O | 100 µg/mL in water/diethylamine (80/20 v/v) | 1.2 mL |
| ULM-8323-1.2 | Ammeline (desethyldesisopropylhydroxyatrazine) (unlabeled) | C ₃ H ₅ N ₅ O | 100 µg/mL in water/diethylamine (80/20 v/v) | 1.2 mL |
| CLM-4748-1.2 | 1,6-Anhydro-β-D-glucose (levoglucosan) (¹³ C ₆ , 98%) | *C ₆ H ₁₀ O ₅ | 100 µg/mL in DMSO | 1.2 mL |
| ULM-8000-1.2 | 1,6-Anhydro-β-D-glucose (levoglucosan) (unlabeled) | C ₆ H ₁₀ O ₅ | 100 µg/mL in DMSO | 1.2 mL |
| DLM-119-1.2 | (+/-)-Chloramphenicol (ring-D ₄ , benzyl-D ₁ , 98%) | NO ₂ C ₆ D ₄ C ₅ D ₁ H ₅ O ₃ NCl ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-6687-1.2 | (+/-)-Chloramphenicol (unlabeled) | NO ₂ C ₆ H ₄ C ₅ H ₆ O ₃ NCl ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-4633-1.2 | 3-Chloro-1,2-propanediol (~10% 2-chloro-1,3-propanediol)(propane-D ₅ , 98%) | CICD ₂ CDOHCD ₂ OH | 1 mg/mL in methanol | 1.2 mL |
| ULM-7998-1.2 | 3-Chloro-1,2-propanediol (unlabeled) | CICH ₂ CHOHCH ₂ OH | 1 mg/mL in methanol | 1.2 mL |
| CNLM-4661-1.2 | Cyanuric acid (¹³ C ₃ , 99%; ¹⁵ N ₃ , 98%+) CP 90%+ | *C ₃ H ₃ *N ₃ O ₃ | 100 µg/mL in water | 1.2 mL |
| CNLM-4661-10X-1.2 | | | 1000 µg/mL in water | 1.2 mL |
| ULM-8157-1.2 | Cyanuric acid (unlabeled) | C ₃ H ₃ N ₃ O ₃ | 100 µg/mL in water | 1.2 mL |
| NEW DLM-2943-1.2 | 2,6-Di(tert-butyl)-4-methyl phenol (BHT) (D ₂₁ , 98%) | C ₁₅ H ₃ D ₂₁ O | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-7494-1.2 | 2,6-Di(tert-butyl)-4-methyl phenol (BHT) (unlabeled) | C ₁₅ H ₂₄ O | 100 µg/mL in nonane | 1.2 mL |
| DLM-1632-1.2 | Diethylene glycol (D ₈ , 98%) | C ₄ D ₈ H ₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8235-1.2 | Diethylene glycol (unlabeled) | C ₄ H ₁₀ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| CNLM-8150-1.2 | Melamine (¹³ C ₃ , 99%; amino- ¹⁵ N ₃ , 98%) | *C ₃ H ₆ *N ₃ N ₃ | 100 µg/mL in water | 1.2 mL |
| CNLM-8150-10X-1.2 | | | 1000 µg/mL in water | 1.2 mL |
| ULM-8156-1.2 | Melamine (unlabeled) | C ₃ H ₆ N ₃ N ₃ | 100 µg/mL in water | 1.2 mL |
| DLM-4412-25 | (-)Menthol (1,2,6,6-D ₄ , 98%) | C ₁₀ H ₁₆ D ₄ O | neat | 25 mg |
| DLM-4766-1.2 | 2-Methylisoborneol (2-methyl-D ₃ , 98%) | C ₁₁ H ₁₇ D ₃ O | 100 µg/mL in nonane | 1.2 mL |
| CDLM-7279-S | N-Nitrosodimethylamine (¹³ C ₂ , 99%; D ₆ , 98%) | *C ₂ D ₆ N ₂ O | 1 mg/mL in MeCl-D ₂ | 1 mL |
| OLM-7310-1.2 | Perchloric acid, sodium salt (¹⁸ O ₄ , 90%+) | NaCl ¹⁸ O ₄ | 100 µg/mL in water | 1.2 mL |
| ULM-7312-1.2 | Perchloric acid, sodium salt (unlabeled) | NaClO ₄ | 100 µg/ml in water | 1.2 mL |
| CLM-3733-1.2 | o-Phenylphenol (phenyl- ¹³ C ₆ , 99%) | *C ₆ H ₅ C ₆ H ₄ OH | 100 µg/ml in nonane | 1.2 mL |
| ULM-7396-1.2 | o-Phenylphenol (unlabeled) | C ₆ H ₅ C ₆ H ₄ OH | 100 µg/ml in nonane | 1.2 mL |
| CLM-3748-1.2 | p-Phenylphenol (phenyl- ¹³ C ₆ , 99%) CP 96% | *C ₆ H ₅ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW OLM-8283-18O-1.2 | Potassium bromate (¹⁸ O ₃ , 98%) CP 90-95% | KBr ¹⁸ O ₃ | 100 µg/mL in ¹⁸ O water | 1.2 mL |
| ULM-8451-1.2 | Potassium bromate (unlabeled) | KBrO ₃ | 100 µg/mL in water | 1.2 mL |
| CNLM-7221-1.2 | Semicarbazide hydrochloride (SEM) (¹³ C, 99%; ¹⁵ N ₂ , 98%) | *CH ₃ *N ₂ NO·HCl | 100 µg/mL in methanol | 1.2 mL |
| ULM-7187-1.2 | Semicarbazide hydrochloride (SEM) (unlabeled) | CH ₅ N ₃ O· HCl | 100 µg/mL in methanol | 1.2 mL |
| DLM-6083-1.2 | 2,4,6-Trichloroanisole (D ₅ , 98%) | C ₆ D ₂ Cl ₃ OCD ₃ | 1 mg/mL in methanol-D | 1.2 mL |
| ULM-7999-1.2 | 2,4,6-Trichloroanisole (unlabeled) | C ₆ H ₂ Cl ₃ OCH ₃ | 1 mg/mL in methanol | 1.2 mL |
| DLM-2080-1.2 | 1,2,3-Trichloropropane (D ₅ , 98%) CP 95% | CD ₂ ClCDCICD ₂ Cl | 1 mg/mL in methanol | 1.2 mL |
| ULM-6911-1.2 | 1,2,3-Trichloropropane (unlabeled) | CH ₂ ClCHClCH ₂ Cl | 1 mg/mL in methanol | 1.2 mL |
| NEW DLM-4444-0.1 | Urethane (ethyl carbamate) (ethyl-D ₅ , 98%) | C ₃ H ₂ D ₅ NO ₂ | neat | 0.1 g |

Please also see the sections on PCBs, pesticides, PAHs and priority pollutants for other products that can be used in food and water analysis.

Phthalate and Phthalate Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|----------------------------|--|---|----------------------|---------|
| DLM-1369-1.2 | Benzyl butyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ [CO ₂ (CH ₂) ₃ CH ₃][CH ₂ C ₆ H ₅] | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1369-0.1 | | | neat | 0.1 g |
| ULM-7551-1.2 | Benzyl butyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₃ CH ₃][CH ₂ C ₆ H ₅] | 100 µg/mL in nonane | 1.2 mL |
| CLM-4675-1.2 | Bis(2-ethylhexyl) adipate (adipate- ¹³ C ₆ , 99%) | (*CH ₂) ₄ [*CO ₂ [CH ₂ CH(C ₂ H ₅)C ₄ H ₉]] ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6566-1.2 | Bis(2-ethylhexyl) adipate (unlabeled) | (CH ₂) ₄ [CO ₂ [CH ₂ CH(C ₂ H ₅)C ₄ H ₉]] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-6238 | Bis(2-ethylhexyl) phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | C ₂₀ *C ₄ H ₃₈ O ₄ | | Inquire |
| DLM-1368-1.2 | Bis(2-ethylhexyl) phthalate (ring-D ₄ , 98%) | C ₆ D ₄ [CO ₂ CH ₂ CH(C ₂ H ₅)C ₄ H ₉] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1368-0.1 | | | neat | 0.1 g |
| NEW DLM-1368-0.25 | | | neat | 0.25 g |
| ULM-6241-1.2 | Bis(2-ethylhexyl) phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₂ CH ₃ (CH ₂) ₃ CH ₃)] ₂ | 1000 µg/mL in nonane | 1.2 mL |
| CLM-4670-1.2 | Dicyclohexyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ (*CO ₂ C ₆ H ₁₁) ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-8785-1.2 | Dicyclohexyl phthalate (unlabeled) | C ₆ H ₄ (CO ₂ C ₆ H ₁₁) ₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1629-1.2 | Diethyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ (CO ₂ CH ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1629-0.1 | | | neat | 0.1 g |
| NEW DLM-1629-0.25 | | | neat | 0.25 g |
| ULM-6174-1.2 | Diethyl phthalate (unlabeled) | C ₆ H ₄ (CO ₂ CH ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9767 | Diisononyl phthalate (unlabeled) | C ₂₆ H ₄₂ O ₄ | | Inquire |
| DLM-1366-1.2 | Dimethyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ (CO ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1366-0.1 | | | neat | 0.1 g |
| NEW ULM-6783-1.2 | Dimethyl phthalate (unlabeled) | C ₆ H ₄ (CO ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1367-1.2 | Di-n-butyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ [CO ₂ (CH ₂) ₃ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1367-0.1 | | | neat | 0.1 g |
| NEW DLM-1367-0.25 | | | neat | 0.25 g |
| NEW ULM-7466-1.2 | Di-n-butyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₃ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4669-1.2 | Di-n-hexyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₅ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7434-1.2 | Di-n-hexyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₅ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1630-1.2 | Di-n-octyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ [CO ₂ (CH ₂) ₇ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1630-0.1 | | | neat | 0.1 g |
| ULM-6129-1.2 | Di-n-octyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₇ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4668-1.2 | Di-n-pentyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₄ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7433-1.2 | Di-n-pentyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₄ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4671 | Di-n-propyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | C ₂ C ₄ H ₄ -1,2-[*CO ₂ (CH ₂) ₂ CH ₃] ₂ | | Inquire |
| CLM-4591-MT-1.2 | Monobenzyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ C ₆ H ₅][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| ULM-6149-MT-1.2 | Monobenzyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ C ₆ H ₅][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4590-MT-1.2 | Mono-n-butyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₃ CH ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-6148-MT-1.2 | Mono-n-butyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₃ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-8232-MT-1.2 | Mono-[2-(carboxymethyl) hexyl] phthalate (DEHP Metabolite IV) (¹³ C ₄ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ (CH ₂) ₅ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-8233-MT-1.2 | Mono-[2-(carboxymethyl) hexyl] phthalate (DEHP Metabolite IV) (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ (CH ₂) ₅ CH ₃ CO ₂][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-6847-MT-1.2 | Mono-(3-carboxypropyl) phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₃ CO ₂ H][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-6848-MT-1.2 | Mono-(3-carboxypropyl) phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₃ CO ₂ H][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4592-MT-1.2 | Monocyclohexyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ C ₆ H ₁₁][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-7394-MT-1.2 | Monocyclohexyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ C ₆ H ₁₁][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |

Phthalate and Phthalate Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|----------------------------|--|--|-------------------|--------|
| NEW CLM-4584-MT-1.2 | Mono-2-ethylhexyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂) ₃ CH ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4583-MT-1.2 | Mono-2-ethylhexyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂) ₃ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-8148-MT-1.2 | Mono-(2-ethyl-5-carboxypentyl) phthalate (DEHP Metabolite V) (¹³ C ₄ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ (CH ₂ CH ₃)(CH ₂) ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-8149-MT-1.2 | Mono-(2-ethyl-5-carboxypentyl) phthalate (DEHP Metabolite V) (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ (CH ₂ CH ₃)(CH ₂) ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-6641-MT-1.2 | Mono-(2-ethyl-5-hydroxyhexyl)phthalate (DEHP Metabolite IX) (¹³ C ₄ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂ CH ₂ CH(OH)CH ₃)[*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4662-MT-1.2 | Mono-(2-ethyl-5-hydroxyhexyl)phthalate (DEHP Metabolite IX) (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂ CH ₂ CH(OH)CH ₃)[CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-6640-MT-1.2 | Mono-(2-ethyl-5-oxohexyl)phthalate (DEHP Metabolite VI) (¹³ C ₄ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂ CH ₂ COCH ₃)[*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4663-MT-1.2 | Mono-(2-ethyl-5-oxohexyl)phthalate (DEHP Metabolite VI) (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₂ CH ₃)CH ₂ CH ₂ COCH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4586-MT-1.2 | Monoethyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₂ CH ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4585-MT-1.2 | Monoethyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4820 | Mono-3-hydroxybutyl phthalate (unlabeled) | C ₁₂ H ₁₄ O ₅ | Inquire | |
| NEW ULM-7919-MT-1.2 | Monoisobutyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₃) ₂][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4588 | Monoisodecyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | C ₁₄ *C ₄ H ₂₆ O ₄ | Inquire | |
| NEW ULM-4652-MT-1.2 | Monoisodecyl phthalate (mono-3,7-dimethyloctyl phthalate) (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₂ CH(CH ₃ (CH ₂) ₃ CH ₃) ₂][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4587-MT-1.2 | Monoisononyl phthalate (mono-3,5,5-trimethylhexyl phthalate) (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₆ CH(CH ₃) ₂][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4651-MT-1.2 | Monoisononyl phthalate (mono-3,5,5-trimethylhexyl phthalate) (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₆ CH(CH ₃) ₂][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-7395-MT-1.2 | Monoisopropyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH(CH ₃) ₂][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4594 | Mono-2-methoxyethyl phthalate (unlabeled) | C ₁₁ H ₁₂ O ₅ | Inquire | |
| NEW CLM-6071-MT-1.2 | Monomethyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ CH ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-6697-MT-1.2 | Monomethyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-6225 | Monomethyl isophthalate (ring- ¹³ C ₆ , 99%) | *C ₆ C ₃ H ₈ O ₄ | Inquire | |
| NEW ULM-6226 | Monomethyl isophthalate (unlabeled) | C ₉ H ₈ O ₄ | Inquire | |
| NEW CLM-4589-MT-1.2 | Mono-n-octyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₇ CH ₃][*CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-4593-MT-1.2 | Mono-n-octyl phthalate (unlabeled) | C ₂ C ₄ H ₄ [CO ₂ (CH ₂) ₇ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-7393-MT-1.2 | Mono-n-pentyl phthalate (unlabeled) | C ₂ C ₄ H ₄ [CO ₂ (CH ₂) ₄ CH ₃][CO ₂ H] | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-4323-MT-1.2 | Phthalic acid (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ *(COOH) ₂ | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-8301-MT-1.2 | Phthalic acid (unlabeled) | C ₆ H ₄ (COOH) ₂ | 100 µg/mL in MTBE | 1.2 mL |

Bisphenol Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|---------------------------|--------|
| CLM-4325-1.2 | Bisphenol A (ring- ¹³ C ₁₂ , 99%) | ([*] C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7106-1.2 | Bisphenol A (unlabeled) | (C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8654-1.2 | 2,4'-Bisphenol A (unlabeled) | (C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW DLM-9193-1.2 | Bisphenol A diglycidyl ether (BADGE) (diglycidyl-D ₁₀ , 98%) | C ₂₁ H ₁₄ D ₁₀ O ₄ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9857-1.2 | Bisphenol A diglycidyl ether (BADGE) (unlabeled) | C ₂₁ H ₁₄ H ₁₀ O ₄ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9831-1.2 | Bisphenol A β-D-glucuronide (unlabeled) CP 90% | C ₂₁ H ₂₄ O ₈ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9832-1.2 | Bisphenol A bis-(β-D-glucoronide) disodium salt (unlabeled) CP 90% | C ₂₇ H ₃₀ Na ₂ O ₁₄ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9833-1.2 | Bisphenol A bisulfate disodium salt (unlabeled) CP 90% | C ₁₅ H ₁₄ Na ₂ O ₈ S ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-9776-1.2 | Bisphenol AF (ring- ¹³ C ₁₂ , 99%) | ([*] C ₆ H ₄ OH) ₂ C(CF ₃) ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9779-1.2 | Bisphenol AF (unlabeled) | (C ₆ H ₄ OH) ₂ C(CF ₃) ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9830-1.2 | Bisphenol AP (unlabeled) | CH ₃ C(C ₆ H ₅)(C ₆ H ₄ OH) ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-9851-1.2 | Bisphenol B (ring- ¹³ C ₁₂ , 99%) | C ₄ *C ₁₂ H ₁₈ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9852-1.2 | Bisphenol B (unlabeled) | C ₁₆ H ₁₈ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9826-1.2 | Bisphenol E (unlabeled) | CH ₃ CH(C ₆ H ₄ O ₄) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-9866-1.2 | Bisphenol F (ring- ¹³ C ₁₂ , 99%) | CH ₂ (*C ₆ H ₄ OH) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9827-1.2 | Bisphenol F (unlabeled) | (C ₆ H ₄ OH) ₂ CH ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-9867-1.2 | Bisphenol F diglycidyl ether (BFDGE) (ring- ¹³ C ₁₂ , 99%) | C ₇ *C ₁₂ H ₂₀ O ₄ | Inquire | |
| NEW ULM-9868-1.2 | Bisphenol F diglycidyl ether (BFDGE) (unlabeled) | C ₁₉ H ₂₀ O ₄ | Inquire | |
| NEW ULM-9829-1.2 | Bisphenol P (unlabeled) | C ₆ H ₄ [C(CH ₃) ₂ C ₆ H ₄ OH] ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-9319-1.2 | Bisphenol S (¹³ C ₁₂ , 98%) | *C ₁₂ H ₁₀ O ₄ S | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9320-1.2 | Bisphenol S (unlabeled) | C ₁₂ H ₁₀ O ₄ S | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9828-1.2 | Bisphenol Z (unlabeled) | C ₆ H ₁₀ (C ₆ H ₄ OH) ₂ | 100 µg/mL in methanol | 1.2 mL |

Perfluorinated Compound Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|----------------------|--------|
| CLM-8505-1.2 | Perfluorooctanesulfonate (PFOS), sodium salt (¹³ C ₈ , 99%) | *C ₈ F ₁₇ NaO ₃ S | 50 µg/mL in methanol | 1.2 mL |
| ULM-9001-1.2 | Perfluorooctanesulfonate (PFOS), sodium salt (unlabeled) | C ₈ F ₁₇ NaO ₃ S | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-8097-1.2 | Perfluorobutyric acid (PFBA) (unlabeled) | CF ₃ CF ₂ CF ₂ COOH | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-9515-1.2 | Perfluoropentanoic acid (PPFA) (unlabeled) | CF ₃ (CF ₂) ₃ COOH | 50 µg/mL in methanol | 1.2 mL |
| CLM-8340-1.2 | Perfluorohexanoic acid (PFHxA), sodium salt (¹³ C ₆ , 99%) | *CF ₃ (*CF ₂) ₄ CO ₂ Na | 50 µg/mL in methanol | 1.2 mL |
| ULM-8342-1.2 | Perfluorohexanoic acid (PFHxA), sodium salt (unlabeled) | CF ₃ (CF ₂) ₄ CO ₂ Na | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-9516-1.2 | Perfluoroheptanoic acid (PFHpA) (unlabeled) | CF ₃ (CF ₂) ₅ CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| CLM-8005-1.2 | Perfluorooctanoic acid (PFOA) (¹³ C ₈ , 99%) | *CF ₃ (*CF ₂) ₆ *CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| ULM-7451-1.2 | Perfluorooctanoic acid (PFOA) (unlabeled) | CF ₃ (CF ₂) ₆ CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| CLM-8060-1.2 | Perfluorononanoic acid (PFNA) (¹³ C ₉ , 99%) | *CF ₃ (*CF ₂) ₇ *CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| ULM-8066-1.2 | Perfluorononanoic acid (PFNA) (unlabeled) | CF ₃ (CF ₂) ₇ CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| CLM-8172-1.2 | Perfluorodecanoic acid (PFDA) (¹³ C ₉ , 99%) | CF ₃ (*CF ₂) ₈ *CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| ULM-8067-1.2 | Perfluorodecanoic acid (PFDA) (unlabeled) | CF ₃ (CF ₂) ₈ CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| NEW CLM-8789-1.2 | Perfluoroundecanoic acid (PFUA), sodium salt (¹³ C ₉ , 99%) | CF ₃ (*CF ₂) ₉ CO ₂ Na | 50 µg/mL in methanol | 1.2 mL |
| ULM-8084-1.2 | Perfluoroundecanoic acid (PFUA), sodium salt (unlabeled) | CF ₃ (CF ₂) ₉ CO ₂ Na | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-8068-1.2 | Perfluorododecanoic acid (PFDoA) (unlabeled) | CF ₃ (CF ₂) ₁₀ CO ₂ H | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-9955-1.2 | Perfluorotridecanoic acid (PTrDA) (unlabeled) | CF ₃ (CF ₂) ₁₁ CO ₂ H | Inquire | |
| NEW ULM-9956-1.2 | Perfluorotetradecanoic acid (PFTeDA) (unlabeled) | CF ₃ (CF ₂) ₁₂ CO ₂ H | Inquire | |

Nonylphenol, Nonylphenol Ethoxylate, and Nonylphenol Carboxylate Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------|--|---|---------------------------|----------------------|
| CLM-8356-1.2 | 4-(1,3-Dimethyl-1-ethylpentyl) phenol (ring- ¹³ C ₆ , 99%) | (CH ₃ CH ₂ CH)(CH ₃)CH ₂ (CH ₃) (CH ₂ CH ₃)C* ₆ H ₄ OH | | Inquire |
| NEW ULM-8360-1.2 | 4-(1,3-Dimethyl-1-ethylpentyl) phenol (unlabeled) | (CH ₃ CH ₂ CH)(CH ₃)CH ₂ (CH ₃) (CH ₂ CH ₃)CC ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-8357-1.2 | 4-(1,4-Dimethyl-1-ethylpentyl) phenol (ring- ¹³ C ₆ , 99%) | (CH ₃) ₂ C(CH ₂) ₂ (CH ₃)(CH ₂ CH ₃) C* ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8361-1.2 | 4-(1,4-Dimethyl-1-ethylpentyl) phenol (unlabeled) | (CH ₃) ₂ C(CH ₂) ₂ (CH ₃)(CH ₂ CH ₃) CC ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-8359-1.2 | 4-(1-Ethyl-1-methylhexyl) phenol (ring- ¹³ C ₆ , 99%) | [(CH ₃)(CH ₂) ₄](CH ₃)(CH ₂ CH ₃) C* ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8363-1.2 | 4-(1-Ethyl-1-methylhexyl) phenol (unlabeled) | [(CH ₃)(CH ₂) ₄](CH ₃)(CH ₂ CH ₃) CC ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| CLM-8358-1.2 | 4-(1,1,5-Trimethylhexyl) phenol (ring- ¹³ C ₆ , 99%) | (CH ₃) ₂ C(CH ₂) ₃ (CH ₃) ₂ (CH ₂ CH ₃)C* ₆ H ₄ OH | | Inquire |
| NEW ULM-8362-1.2 | 4-(1,1,5-Trimethylhexyl) phenol (unlabeled) | (CH ₃) ₂ C(CH ₂) ₃ (CH ₃) ₂ (CH ₂ CH ₃)CC ₆ H ₄ OH | 100 µg/mL in methanol | 1.2 mL |
| CLM-4306-1.2 | p-n-Nonylphenol (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ * ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4306-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4559-1.2 | p-n-Nonylphenol (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-4559-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| CLM-4512-1.2 | p-n-Nonylphenol monoethoxylate (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ * ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4512-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4520-1.2 | p-n-Nonylphenol monoethoxylate (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-4520-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4520-SA-5X-1.2 | | | 500 µg/mL in acetonitrile | 1.2 mL |
| CLM-4307-1.2 | p-n-Nonylphenol diethoxylate (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ * ₆ H ₄ O(CH ₂) ₂ O (CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4307-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4521-1.2 | p-n-Nonylphenol diethoxylate (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ O(CH ₂) ₂ O (CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-4521-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4521-SA-5X-1.2 | | | 500 µg/mL in acetonitrile | 1.2 mL |
| CLM-4516-1.2 | p-n-Nonylphenol triethoxylate (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ * ₆ H ₄ O(CH ₂) ₂ O (CH ₂) ₂ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| CP 90% | | | | |
| ES-4157 | p-n-Nonylphenol + mono-/di-/tri-ethoxylates (set of individual standards) 1 ampoule each of CLM-4306-1.2, CLM-4512-1.2, CLM-4307-1.2 and CLM-4516-1.2 | | | Set of 4 × 1.2 mL |
| ULM-6560-1.2 | p-Nonylphenol – technical grade (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-7146-1.2 | Nonylphenol monoethoxylate-branched isomers (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-7147-1.2 | Nonylphenol diethoxylate-branched isomers (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ (OCH ₂ CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-4688-1.2 | Nonylphenoxyacetic acid – ring/chain isomers (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ OCH ₂ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |
| ULM-4690-1.2 | p-n-Nonylphenoxyethoxyacetic acid (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ O(CH ₂) ₂ OCH ₂ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |

Chlorinated Paraffin Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------|---|--|---------------------|--------|
| NEW CLM-9000-1.2 | 1,5,5,6,6,10-Hexachlorodecane (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8917-1.2 | 1,5,5,6,6,10-Hexachlorodecane (unlabeled) | C ₁₀ H ₁₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9679-1.2 | 1,1,1,3,10,12,12,12-Octachlorododecane (¹³ C ₁₂ , 99%) | *C ₁₂ H ₁₈ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9485-1.2 | 1,1,1,3,10,12,12,12-Octachlorododecane (unlabeled) | C ₁₂ H ₁₈ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |

Nitrosamine Standards

| Catalog No. | Compound | Formula | Concentration | Amount | |
|--------------|---|--|--------------------------------|--------------------------------|--------|
| DLM-7779-S | <i>N</i> -Nitrodimethylamine (D_6 , 98%) | $C_2D_6N_2O_2$ | 1 mg/mL in MeCl-D ₂ | 1 mL | |
| ULM-7780-S | <i>N</i> -Nitrodimethylamine (unlabeled) | $C_2H_6N_2O_2$ | 1 mg/mL in MeCl | 1 mL | |
| DLM-7982-S | <i>N</i> -Nitrosodiethylamine (D_{10} , 98%) | $(C_2D_5)_2NNO$ | 1 mg/mL in MeCl-D ₂ | 1 mL | |
| ULM-7984-1.2 | <i>N</i> -Nitrosodiethylamine (unlabeled) | $(C_2H_5)_2NNO$ | 1 mg/mL in MeCl | 1.2 mL | |
| CDLM-7279-S | <i>N</i> -Nitrosodimethylamine (¹³ C ₂ , 99%; D_6 , 98%) | * $C_2D_6N_2O$ | 1 mg/mL in MeCl-D ₂ | 1 mL | |
| DLM-2130-S | <i>N</i> -Nitrosodimethylamine (D_6 , 98%) | $C_2D_6N_2O$ | 1 mg/mL in MeCl-D ₂ | 1 mL | |
| NLM-7647-S | <i>N</i> -Nitrosodimethylamine (¹⁵ N ₂ , 98%) | $(C_2H_5)_2^*N^*NO$ | 1 mg/mL in MeCl | 1 mL | |
| NEW | ULM-9042-S | <i>N</i> -Nitrosodimethylamine (unlabeled) | $(C_2H_5)_2NNO$ | 1 mg/mL in MeCl | 1 mL |
| | DLM-3098-S | <i>N</i> -Nitrosodiphenylamine (2,2',4,4',6,6'-D ₆ , 98%) | $(C_6D_3H_2)_2NN=O$ | 1 mg/mL in MeCl-D ₂ | 1 mL |
| | ULM-7219-1.2 | <i>N</i> -Nitrosodiphenylamine (unlabeled) | $C_{12}H_{10}N_2O$ | 1 mg/mL in MeCl | 1.2 mL |
| | DLM-2131-S | <i>N</i> -Nitrosodi- <i>n</i> -propylamine (D_{14} , 98%) | $C_6D_{14}N_2O$ | 1 mg/mL in MeCl-D ₂ | 1 mL |
| | ULM-6637-S | <i>N</i> -Nitrosodi- <i>n</i> -propylamine (unlabeled) | $C_6H_{14}N_2O$ | 1 mg/mL in MeCl | 1 mL |
| | DLM-8254-1.2 | <i>N</i> -Nitrosomorpholine (D_8 , 98%) | $CD_8N_2O_2$ | 1 mg/mL in MeCl-D ₂ | 1.2 mL |
| | ULM-8255-1.2 | <i>N</i> -Nitrosomorpholine (unlabeled) CP 96% | $CH_8N_2O_2$ | 1 mg/mL in MeCl | 1.2 mL |
| | DLM-8252-1.2 | <i>N</i> -Nitrosopyrrolidine (D_8 , 98%) | $C_4D_8N_2O$ | 1 mg/mL in MeCl-D ₂ | 1.2 mL |
| | ULM-8253-1.2 | <i>N</i> -Nitrosopyrrolidine (unlabeled) | $C_4H_8N_2O$ | 1 mg/mL in MeCl | 1.2 mL |

Halogenated and Substituted Benzene, Phenol, and Anisole Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|-----------------------|--------|
| CLM-871-0.5 | Bromobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_5\text{Br}$ | neat | 0.5 g |
| DLM-398-5 | Bromobenzene (D ₅ , 99%) | C ₆ D ₅ Br | neat | 5 g |
| DLM-398-10 | | | neat | 10 g |
| DLM-398-25 | | | neat | 25 g |
| CLM-2268-1.2 | 4-Bromophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{BrOH}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-6917-1.2 | 4-Bromophenol (unlabeled) | C ₆ H ₄ BrOH | 100 µg/mL in toluene | 1.2 mL |
| DLM-263-1 | Chlorobenzene (D ₅ , 99%) | C ₆ D ₅ Cl | neat | 1 g |
| DLM-263-5 | | | neat | 5 g |
| DLM-1638-0.1 | 2-Chlorophenol (ring-D ₄ , 99%) | C ₆ D ₄ ClOH | neat | 0.1 g |
| DLM-1638-0.25 | | | neat | 0.25 g |
| CLM-1913-1.2 | 4-Chlorophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{ClOH}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-7420-1.2 | 4-Chlorophenol (unlabeled) | C ₆ H ₄ ClOH | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9373-1.2 | 2,4-Dibromoanisole (ring- $^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{Br}_2\text{H}_3\text{OCH}_3$ | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-9369-1.2 | 2,4-Dibromoanisole (unlabeled) | C ₆ Br ₂ H ₃ OCH ₃ | 100 µg/mL in toluene | 1.2 mL |
| CLM-1340-0.1 | 1,4-Dibromobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{Br}_2$ | neat | 0.1 g |
| DLM-341-5 | 1,4-Dibromobenzene (D ₄ , 98%) | C ₆ D ₄ Br ₂ | neat | 5 g |
| CLM-6058-1.2 | 2,4-Dibromophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_3\text{Br}_2\text{OH}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-6918-1.2 | 2,4-Dibromophenol (unlabeled) | C ₆ H ₃ Br ₂ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-8007-1.2 | 2,6-Dibromophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_3\text{Br}_2\text{OH}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-7603-1.2 | 2,6-Dibromophenol (unlabeled) | C ₆ H ₃ Br ₂ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-126-1.2 | 1,2-Dichlorobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{Cl}_2$ | 100 µg/mL in isoctane | 1.2 mL |
| DLM-158-1 | 1,2-Dichlorobenzene (D ₄ , 99%) | C ₆ D ₄ Cl ₂ | neat | 1 g |
| DLM-158-5 | | | neat | 5 g |
| NEW ULM-7415-1.2 | 1,2-Dichlorobenzene (unlabeled) | C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-4484-1.2 | 1,3-Dichlorobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{Cl}_2$ | 100 µg/mL in isoctane | 1.2 mL |
| NEW DLM-2139-0.1 | 1,3-Dichlorobenzene (D ₄ , 98%) | C ₆ D ₄ Cl ₂ | neat | 0.1 g |
| NEW ULM-7431-1.2 | 1,3-Dichlorobenzene (unlabeled) | C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-1518-1 | 1,4-Dichlorobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{Cl}_2$ | neat | 1 mg |
| DLM-268-5 | 1,4-Dichlorobenzene (D ₄ , 98%) | C ₆ D ₄ Cl ₂ | neat | 5 g |
| DLM-1359-0.1 | 2,4-Dichlorophenol (ring-D ₃ , 98%) | C ₆ D ₃ Cl ₂ OH | neat | 0.1 g |
| DLM-1359-0.5 | | | neat | 0.5 g |
| DLM-1669-0.1 | 2,4-Dichlorophenol (D ₄ , 98%) | C ₆ D ₃ Cl ₂ OD | neat | 0.1 g |
| ULM-6822-1.2 | 2,4-Dichlorophenol (unlabeled) | C ₆ H ₃ Cl ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-1365-1.2 | 2,5-Dichlorophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_3\text{Cl}_2\text{OH}$ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9066-1.2 | 2,5-Dichlorophenol (unlabeled) | C ₆ H ₃ Cl ₂ OH | 100 µg/mL in methanol | 1.2 mL |
| CLM-1921-1.2 | Hexabromobenzene ($^{13}\text{C}_6$, 99%) | * C_6Br_6 | 100 µg/mL in toluene | 1.2 mL |
| ULM-7607-1.2 | Hexabromobenzene (unlabeled) | C ₆ Br ₆ | 100 µg/mL in toluene | 1.2 mL |
| CLM-351-1.2 | Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | * C_6Cl_6 | 100 µg/mL in nonane | 1.2 mL |
| CLM-351-0.1 | | | neat | 0.1 g |
| ULM-6130-1.2 | Hexachlorobenzene (unlabeled) | C ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-8992-1.2 | Pentabromoanisole ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{CH}_3\text{Br}_5\text{O}$ | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-8991-1.2 | Pentabromoanisole (unlabeled) | C ₇ H ₃ Br ₅ O | 100 µg/mL in toluene | 1.2 mL |
| CLM-1959-1.2 | Pentabromophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{Br}_5\text{OH}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-6922-1.2 | Pentabromophenol (unlabeled) | C ₆ Br ₅ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-8003-1.2 | Pentachloroanisole ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{CH}_3\text{Cl}_5\text{O}$ | 100 µg/mL in toluene | 1.2 mL |
| ULM-7605-1.2 | Pentachloroanisole (unlabeled) | C ₆ CH ₃ Cl ₅ O | 100 µg/mL in toluene | 1.2 mL |
| CLM-2050-1.2 | Pentachlorobenzene ($^{13}\text{C}_6$, 99%) | * C_6HCl_5 | 100 µg/mL in isoctane | 1.2 mL |
| ULM-7234-1.2 | Pentachlorobenzene (unlabeled) | C ₆ HCl ₅ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-1955-1.2 | Pentachloronitrobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{Cl}_5\text{NO}_2$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7597-1.2 | Pentachloronitrobenzene (unlabeled) | C ₆ Cl ₅ NO ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-661-1.2 | Pentachlorophenol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{Cl}_5\text{OH}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-661-0.01 | | | neat | 0.01 g |
| ULM-6894-1.2 | Pentachlorophenol (unlabeled) | C ₆ Cl ₅ OH | 100 µg/mL in nonane | 1.2 mL |

(continued on next page)

Halogenated and Substituted Benzene, Phenol, and Anisole Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|----------------------------|--|--|-----------------------|---------|
| CLM-1996-1.2 | 2,3,4,5-Tetrabromophenol (¹³ C ₆ , 99%) | *C ₆ HBr ₄ OH | 100 µg/mL in toluene | 1.2 mL |
| ULM-6778-1.2 | 2,3,4,5-Tetrabromophenol (unlabeled) | C ₆ HBr ₄ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-1982-1.2 | 1,2,3,4-Tetrachlorobenzene (¹³ C ₆ , 99%) | *C ₆ H ₂ Cl ₄ | 100 µg/mL in isoctane | 1.2 mL |
| ULM-6195-1.2 | 1,2,3,4-Tetrachlorobenzene (unlabeled) | C ₆ H ₂ Cl ₄ | 100 µg/mL in isoctane | 1.2 mL |
| ULM-7599-1.2 | 1,2,3,5-Tetrachlorobenzene (unlabeled) | C ₆ H ₂ Cl ₄ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-585-0.1 | 1,2,4,5-Tetrachlorobenzene (¹³ C ₆ , 99%) | *C ₆ H ₂ Cl ₄ | neat | 0.1 g |
| CLM-585-5 | | | neat | 5 mg |
| DLM-1177-1 | 1,2,4,5-Tetrachlorobenzene (D ₂ , 98%) | C ₆ D ₂ Cl ₄ | neat | 1 g |
| DLM-1177-5 | | | neat | 5 g |
| ULM-7598-1.2 | 1,2,4,5-Tetrachlorobenzene (unlabeled) | C ₆ H ₂ Cl ₄ | 100 µg/mL in isoctane | 1.2 mL |
| ULM-2428-0.1 | 2,3,4,5-Tetrachlorophenol (unlabeled) | C ₆ H ₂ Cl ₄ O | neat | 0.1 g |
| ULM-2429-0.1 | 2,3,4,6-Tetrachlorophenol (unlabeled) | C ₆ H ₂ Cl ₄ O | neat | 0.1 g |
| ULM-2430-0.1 | 2,3,5,6-Tetrachlorophenol (unlabeled) | C ₆ H ₂ Cl ₄ O | neat | 0.1 g |
| NEW CLM-9372-1.2 | 2,4,5-Tribromoanisole (ring- ¹³ C ₆ , 99%) | *C ₆ CH ₅ Br ₃ O | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-9367-1.2 | 2,4,5-Tribromoanisole (unlabeled) | C ₇ H ₅ Br ₃ O | 100 µg/mL in toluene | 1.2 mL |
| NEW CLM-6744-1.2 | 2,4,6-Tribromoanisole (ring- ¹³ C ₆ , 99%) | *C ₆ CH ₅ Br ₃ O | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-9370-1.2 | 2,4,6-Tribromoanisole (unlabeled) | C ₇ H ₅ Br ₃ O | 100 µg/mL in toluene | 1.2 mL |
| CLM-7488 | 2,3,4-Tribromophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Br ₃ OH | | Inquire |
| NEW CLM-2235-1.2 | 2,3,5-Tribromophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| NEW ULM-6919-1.2 | 2,3,5-Tribromophenol (unlabeled) | C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-6151-1.2 | 2,4,5-Tribromophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| ULM-6084-1.2 | 2,4,5-Tribromophenol (unlabeled) | C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-6743-1.2 | 2,4,6-Tribromophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| ULM-4210-1.2 | 2,4,6-Tribromophenol (unlabeled) | C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| CLM-1836-1.2 | 3,4,5-Tribromophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Br ₃ OH | 100 µg/mL in toluene | 1.2 mL |
| NEW DLM-9198 | 2,4,6-Trichloroanisole (methyl-D ₃ , 99%) | C ₆ H ₂ Cl ₃ OCD ₃ | | Inquire |
| DLM-6083-1.2 | 2,4,6-Trichloroanisole (D ₅ , 98%) | C ₆ D ₂ Cl ₃ OCD ₃ | 1 mg/mL in methanol-D | 1.2 mL |
| DLM-6083-0.1 | | | neat | 0.1 g |
| ULM-7999-1.2 | 2,4,6-Trichloroanisole (unlabeled) | C ₆ H ₂ Cl ₃ OCH ₃ | 1 mg/mL in methanol | 1.2 mL |
| DLM-1972-0.1 | 1,2,3-Trichlorobenzene (D ₃ , 98%) | C ₆ D ₃ Cl ₃ | neat | 0.1 g |
| DLM-1178-0.1 | 1,2,4-Trichlorobenzene (D ₃ , 98%) | C ₆ D ₃ Cl ₃ | neat | 0.1 g |
| DLM-1178-1 | | | neat | 1 g |
| DLM-1178-5 | | | neat | 5 g |
| DLM-799-1 | 1,3,5-Trichlorobenzene (D ₃ , 98%) | C ₆ D ₃ Cl ₃ | neat | 1 g |
| CLM-513-1 | 2,4,5-Trichlorophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Cl ₃ OH | 100 µg/mL in methanol | 1 mL |
| CLM-513-SI-1.2 | | | 100 µg/mL in isoctane | 1.2 mL |
| DLM-2143-0.1 | 2,4,5-Trichlorophenol (ring-D ₂ , 98%) | C ₆ D ₂ Cl ₃ OH | neat | 0.1 g |
| ULM-7525-1.2 | 2,4,5-Trichlorophenol (unlabeled) | C ₆ H ₂ Cl ₃ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-1804-1.2 | 2,4,6-Trichlorophenol (¹³ C ₆ , 99%) | *C ₆ H ₂ Cl ₃ OH | 100 µg/mL in methanol | 1.2 mL |
| CLM-1804-SI-1.2 | | | 100 µg/mL in isoctane | 1.2 mL |
| DLM-3093-0.01 | 2,4,6-Trichlorophenol (ring-D ₂ , 98%) | C ₆ D ₂ Cl ₃ OH | neat | 0.01 g |
| DLM-3093-0.1 | | | neat | 0.1 g |
| ULM-7600-1.2 | 2,4,6-Trichlorophenol (unlabeled) | C ₆ H ₂ Cl ₃ OH | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-7600-SI-1.2 | | | 100 µg/mL in isoctane | 1.2 mL |

Please also see the priority pollutant mixtures section for halogenated benzene and phenol cocktails.

Endocrine-Disrupting Compounds and Xenoestrogen Standards

| Catalog No. | Compound | Formula | Concentration | Amount | |
|----------------|---|---|---|-------------------------------------|------------|
| CLM-1643-1.2 | Acenaphthene (¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₀ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-108-1.2 | Acenaphthene (D ₁₀ , 98%) | C ₁₂ D ₁₀ | 200 µg/mL in isoctane | 1.2 mL | |
| ULM-7413-1.2 | Acenaphthene (unlabeled) | C ₁₂ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL | |
| CLM-3727-1.2 | Alachlor (ring- ¹³ C ₆ , 99%) CP 96%+ | *C ₆ C ₈ H ₂₀ ClNO ₂ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-4725-1.2 | Aldrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1333-1.2 | Anthracene (¹³ C ₆ , 99%) | *C ₆ C ₈ H ₁₀ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-102-1.2 | Anthracene (D ₁₀ , 98%) | C ₁₄ D ₁₀ | 200 µg/mL in isoctane | 1.2 mL | |
| ULM-7412-1.2 | Anthracene (unlabeled) | C ₁₄ H ₁₀ | 200 µg/mL in isoctane | 1.2 mL | |
| CLM-3737-1.2 | Atrazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₅ H ₁₄ ClN ₅ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-3602-1.2 | Benz[a]anthracene (¹³ C ₆ , 99%) | *C ₆ C ₁₂ H ₁₂ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-610-1.2 | Benz[a]anthracene (D ₁₂ , 98%) | C ₁₈ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL | |
| ULM-2415-I-1.2 | Benz[a]anthracene (unlabeled) | C ₁₈ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL | |
| CLM-2722-1.2 | Benzo[a]pyrene (¹³ C ₄ , 99%) | *C ₄ C ₁₆ H ₁₂ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-258-1.2 | Benzo[a]pyrene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL | |
| NEW | ULM-2412-I-1.2 | Benzo[a]pyrene (unlabeled) | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | CLM-6170-1.2 | Benzo[e]pyrene (¹³ C ₄ , 99%) | *C ₄ C ₁₆ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-257-1.2 | Benzo[e]pyrene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | ULM-7423-1.2 | Benzo[e]pyrene (unlabeled) | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | CLM-3599-1.2 | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | *C ₆ C ₁₄ H ₁ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-2136-1.2 | Benzo[b]fluoranthene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | ULM-2416-I-1.2 | Benzo[b]fluoranthene (unlabeled) | C ₂₀ H ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| NEW | CLM-9590-1.2 | Benzo[j]fluoranthene (¹³ C ₁₂ , 99%) | *C ₁₂ C ₈ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-2411-1.2 | Benzo[j]fluoranthene (unlabeled) | C ₂₀ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | CLM-3756-1.2 | Benzo[k]fluoranthene (¹³ C ₆ , 99%) | *C ₆ C ₁₄ H ₁ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-1923-1.2 | Benzo[k]fluoranthene (D ₁₂ , 98%) | C ₂₀ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| | CLM-9730-1.2 | Benzo[c]phenanthrene (¹³ C ₆ , 99%) | *C ₆ C ₁₂ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-8155-1.2 | Benzo[c]phenanthrene (unlabeled) | C ₁₈ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-183-1.2 | Benzophenone (D ₁₀ , 98%) | C ₆ D ₅ COC ₆ D ₅ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-8303-1.2 | Benzophenone (unlabeled) | (C ₆ H ₅) ₂ CO | 100 µg/mL in nonane | 1.2 mL |
| | DLM-1369-1.2 | Benzyl butyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ [CO ₂ (CH ₂) ₃ CH ₃][CH ₂ C ₆ H ₅] | 100 µg/mL in nonane | 1.2 mL |
| | CLM-2482-1.2 | α-HCH (α-BHC) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| | CLM-3623-1.2 | β-HCH (β-BHC) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 50 µg/mL in nonane | 2 × 1.2 mL |
| | CLM-1282-1.2 | γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| | CLM-4675-1.2 | Bis(2-ethylhexyl) adipate (adipate- ¹³ C ₆ , 99%) | (*CH ₂) ₄ [*CO ₂ [CH ₂ CH(C ₂ H ₅)C ₄ H ₉]] ₂ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-1368-1.2 | Bis(2-ethylhexyl) phthalate (ring-D ₄ , 98%) | C ₆ D ₄ -1,2-[CO ₂ C ₈ H ₁₇] ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-6241-1.2 | Bis(2-ethylhexyl) phthalate (unlabeled) | C ₆ H ₄ [CO ₂ CH ₂ CH(CH ₂ CH ₃)(CH ₂) ₃ CH ₃] ₂ | 1000 µg/mL in nonane | 1.2 mL |
| | CLM-4325-1.2 | Bisphenol A (ring- ¹³ C ₁₂ , 99%) | (*C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| | ULM-7106-1.2 | Bisphenol A (unlabeled) | (C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| | ULM-8654-1.2 | 2,4'-Bisphenol A (unlabeled) | (C ₆ H ₄ OH) ₂ C(CH ₃) ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| | CLM-9319-1.2 | Bisphenol S (¹³ C ₁₂ , 98%) | *C ₁₂ H ₁₀ O ₄ S | 100 µg/mL in methanol | 1.2 mL |
| | ULM-9320-1.2 | Bisphenol S (unlabeled) | C ₁₂ H ₁₀ O ₄ S | 100 µg/mL in methanol | 1.2 mL |
| | CLM-9776-1.2 | Bisphenol AF (ring- ¹³ C ₁₂ , 99%) | *C ₁₂ C ₃ H ₁₀ F ₆ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW | ULM-9779-1.2 | Bisphenol AF (unlabeled) | C ₁₅ H ₁₀ F ₆ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| | CLM-4674-1.2 | n-Butylbenzene (ring- ¹³ C ₆ , 99%) | *C ₆ H ₅ C ₄ H ₉ | 100 µg/mL in nonane | 1.2 mL |
| | CLM-4682-1.2 | Carbaryl (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₁ NO ₂ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-8096-1.2 | Carbaryl (unlabeled) | C ₁₀ H ₇ CO ₂ NHCH ₃ | 100 µg/mL in nonane | 1.2 mL |
| | CLM-1911-1.2 | Carbofuran (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₅ NO ₃ | 100 µg/mL in 1,4-dioxane | 1.2 mL |
| | ULM-7419-1.2 | Carbofuran (unlabeled) | C ₁₂ H ₁₅ NO ₃ | 100 µg/mL in 1,4-dioxane | 1.2 mL |
| | CLM-4792-1.2 | trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | *C ₁₀ H ₈ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW | CLM-4814-1.2 | Chlordecone (kepone) (¹³ C ₁₀ , 99%) | *C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| | ULM-2301-1.2 | Chlordecone (kepone) (unlabeled) | C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| | CLM-4758-1.2 | Chlordene (¹³ C ₁₀ , 99%) | *C ₁₀ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-7443-1.2 | Chlordene (unlabeled) | C ₁₀ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-4360-1.2 | Chlorpyrifos (diethyl-D ₁₀ , 99%) | C ₉ D ₁₀ H ₁ Cl ₃ NO ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| | CLM-3757-1.2 | Chrysene (¹³ C ₆ , 99%) | *C ₆ C ₁₂ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| | DLM-261-1.2 | Chrysene (D ₁₂ , 98%) | C ₁₈ D ₁₂ | 200 µg/mL in toluene-D ₈ | 1.2 mL |
| NEW | ULM-7424-1.2 | Chrysene (unlabeled) | C ₁₈ H ₁₂ | 200 µg/mL in toluene | 1.2 mL |
| | CLM-7293-1.2 | Cyfluthrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |

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Endocrine-Disrupting Compounds and Xenoestrogen Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|---|------------|
| ULM-7454-1.2 | Cyfluthrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7292-1.2 | Cypermethrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7453-1.2 | Cypermethrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| DLM-4461-1.2 | Daidzein (3',5',8-D ₃ , 97%) | C ₁₅ D ₃ H ₇ O ₄ | 60 µg/mL in acetonitrile-D ₃ | 2 × 1.2 mL |
| ULM-4459-1.2 | Daidzein (unlabeled) | C ₁₅ H ₁₀ O ₄ | 60 µg/mL in acetonitrile | 1.2 mL |
| CLM-6999-1.2 | 2,4'-DDD (ring- ¹³ C ₁₂ , 99%) | *C ₁₂ C ₂ H ₁₀ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| DLM-3533-1.2 | 4,4'-DDD (ring-D ₈ , 98%) | C ₁₄ D ₈ H ₄ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4693-1.2 | 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) | (C ₁ *C ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6251-1.2 | 2,4'-DDE (unlabeled) | (ClC ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1627-1.2 | 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | (C ₁ *C ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4692-1.2 | 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | (C ₁ *C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6134-1.2 | 2,4'-DDT (unlabeled) | (ClC ₆ H ₄)CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1281-1.2 | 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | (C ₁ *C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6135-1.2 | 4,4'-DDT (unlabeled) | (ClC ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1148-1.2 | Diazinon (diethyl-D ₁₀ , 98%) | C ₁₂ D ₁₀ H ₁₁ N ₂ O ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| DLM-2943-1.2 | 2,6-Di(tert-butyl)-4-methylphenol (BHT) (D ₂₁ , 98%) | C ₆ D ₂ (CD ₃) ₂ CH ₃ OD | 100 µg/mL in nonane | 1.2 mL |
| CLM-126-1.2 | 1,2-Dichlorobenzene (¹³ C ₆ , 99%) | *C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| NEW ULM-7415-1.2 | 1,2-Dichlorobenzene (unlabeled) | C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-4484-1.2 | 1,3-Dichlorobenzene (¹³ C ₆ , 99%) | *C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| NEW ULM-7431-1.2 | 1,3-Dichlorobenzene (unlabeled) | C ₆ H ₄ Cl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| DLM-1669-0.1 | 2,4-Dichlorophenol (ring-D ₃ , OD, 98%) | C ₆ D ₃ Cl ₂ OD | neat | 0.1 g |
| CLM-1858-1.2 | 2,4-Dichlorophenoxyacetic acid (ring- ¹³ C ₆ , 99%) | Cl ₂ *C ₆ H ₃ OCH ₂ CO ₂ H | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-4726-1.2 | Dieldrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-7230-1.2 | Dieldrin (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| DLM-1629-1.2 | Diethyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ (CO ₂ CH ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6174-1.2 | Diethyl phthalate (unlabeled) | C ₆ H ₄ (CO ₂ CH ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-7151-1.2 | Dimethoate (O,O-dimethyl-D ₆ , 98%) | C ₅ D ₆ H ₆ NO ₃ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-7972-1.2 | Dimethoate (unlabeled) | C ₅ H ₁₂ NO ₃ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-1367-1.2 | Di-n-butyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ (COO(CH ₂) ₃ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-7466-1.2 | Di-n-butyl phthalate (unlabeled) | C ₁₆ H ₂₂ O ₄ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4669-1.2 | Di-n-hexyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ (*CO ₂ (CH ₂) ₅ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7434-1.2 | Di-n-hexyl phthalate (unlabeled) | C ₆ H ₄ (CO ₂ (CH ₂) ₅ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4668-1.2 | Di-n-pentyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ *CO ₂ (CH ₂) ₄ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7433-1.2 | Di-n-pentyl phthalate (unlabeled) | C ₆ H ₄ [CO ₂ (CH ₂) ₄ CH ₃] ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4671 | Di-n-propyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | *C ₂ C ₄ H ₄ [*CO ₂ (CH ₂) ₂ CH ₃] ₂ | Inquire | |
| CLM-6025-1.2 | Endosulfan I (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| DLM-2862-1.2 | Endosulfan I (D ₄ , 97%) | C ₉ D ₄ H ₂ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7447-1.2 | Endosulfan I (unlabeled) | C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-6026-1.2 | Endosulfan II (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7448-1.2 | Endosulfan II (unlabeled) | C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-7531-1.2 | Endosulfan sulfate (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₄ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7990-1.2 | Endosulfan sulfate (unlabeled) | C ₉ H ₆ Cl ₆ O ₄ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-4782-1.2 | Endrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-7444-1.2 | Endrin (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-4815-1.2 | Endrin aldehyde (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4815-50 | | | neat | 50 µg |
| NEW ULM-8958-1.2 | Endrin aldehyde (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4816-1.2 | Endrin ketone (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-4816-50 | | | neat | 50 µg |

Endocrine-Disrupting Compounds and Xenoestrogen Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------------|---|---|---------------------------|--------|
| NEW ULM-8956-1.2 | Endrin ketone (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-3374-1.2 | Epichlorohydrin (¹³ C ₃ , 99%) | *C ₃ H ₅ OCl | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7403-1.2 | Epichlorohydrin (unlabeled) | C ₃ H ₅ OCl | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-4460-1.2 | Genistein (3',5',6,8-D ₄ , 95%) | C ₁₅ D ₄ H ₅ O ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CNLM-4666-1.2 | Glyphosate (2- ¹³ C, 99%; ¹⁵ N, 98%+) CP 96% | HOOC*CH ₂ *NHCH ₂ PO(OH) ₂ | 100 µg/mL in water | 1.2 mL |
| CNLM-4666-10X-1.2 | | | 1000 µg/mL in water | 1.2 mL |
| ULM-6876-1.2 | Glyphosate (unlabeled) | HOOCCH ₂ NHCH ₂ PO(OH) ₂ | 100 µg/mL in water | 1.2 mL |
| CLM-4759-1.2 | Heptachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₇ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-1.2 | Heptachlor (unlabeled) | C ₁₀ H ₅ Cl ₇ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4734-1.2 | cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₇ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2425-1.2 | cis-Heptachlor epoxide (unlabeled) | C ₁₀ H ₅ Cl ₇ O | 100 µg/mL in nonane | 1.2 mL |
| EB-5162 | 2,2',4,4',5,5'-HexaBB (¹³ C ₁₂ , 99%) | *C ₁₂ H ₄ Br ₆ | 40 +/- 4 µg/mL in nonane | 1.2 mL |
| PBB-153-CS | 2,2',4,4',5,5'-HexaBB (Certified Standard) (unlabeled) | C ₁₂ H ₄ Br ₆ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-351-1.2 | Hexachlorobenzene (¹³ C ₆ , 99%) | *C ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6130-1.2 | Hexachlorobenzene (unlabeled) | C ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9000-1.2 | 1,5,5,6,6,10-Hexachlorodecane (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8917-1.2 | 1,5,5,6,6,10-Hexachlorodecane (unlabeled) | C ₁₀ H ₁₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9429-1.2 | Hp-Sed (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9428-1.2 | Hx-Sed (unlabeled) | C ₁₀ H ₁₂ Cl ₆ | 10 µg/mL in nonane | 1.2 mL |
| CLM-3600-1.2 | Indeno[1,2,3-cd]pyrene (¹³ C ₆ , 99%) | *C ₆ C ₁₄ H ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-2148-1.2 | Indeno[1,2,3-cd]pyrene (D ₁₂ , 98%) | C ₂₂ D ₁₂ | 200 µg/mL in isoctane | 1.2 mL |
| CLM-4727-1.2 | Isodrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7442-1.2 | Isodrin (unlabeled) | C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| DLM-4476-1.2 | Malathion (D ₁₀ , 99%) | C ₁₀ D ₁₀ H ₉ O ₆ PS ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-8122-1.2 | Malathion (unlabeled) | C ₁₀ H ₁₉ O ₆ PS ₂ | 100 µg/mL in nonane | 1.2 mL |
| CNLM-8150-1.2 | Melamine (¹³ C ₃ , 99%; amino- ¹⁵ N ₃ , 98%) | *C ₃ H ₆ *N ₃ N ₃ | 100 µg/mL in water | 1.2 mL |
| CNLM-8150-10X-1.2 | | | 1000 µg/mL in water | 1.2 mL |
| ULM-8156-1.2 | Melamine (unlabeled) | C ₃ H ₆ N ₆ | 100 µg/mL in water | 1.2 mL |
| CNLM-7148-1.2 | Methomyl (acetohydroxamate- ¹³ C ₂ , 99%; ¹⁵ N, 98%) | *C ₂ C ₃ H ₁₀ N*NO ₂ S | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8639-1.2 | Methomyl (unlabeled) | C ₅ H ₁₀ N ₂ O ₂ S | 100 µg/mL in methanol | 1.2 mL |
| CLM-4683-1.2 | Methoxychlor (ring- ¹³ C ₁₂ , 99%) | (H ₃ C* _C 6H ₄) ₂ CHCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7440-1.2 | Methoxychlor (unlabeled) | (H ₃ CC ₆ H ₄) ₂ CHCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3712-1.2 | Metolachlor (ring- ¹³ C ₆ , 99%) | *C ₆ C ₉ H ₂₂ CINO ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7314-1.2 | Metolachlor (unlabeled) | C ₁₅ H ₂₂ CINO ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4813-1.2 | Mirex (¹³ C ₁₀ , 99%) | *C ₁₀ Cl ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2427-1.2 | Mirex (unlabeled) | C ₁₀ Cl ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-8246 | Musk ketone (tert-butyl-D ₉ , 98%) | (CD ₃) ₃ CC ₆ (NO ₂) ₂ (CH ₃) ₂ COCH ₃ | Inquire | |
| NEW ULM-8290 | Musk ketone (unlabeled) | (CH ₃) ₃ CC ₆ (NO ₂) ₂ (CH ₃) ₂ COCH ₃ | Inquire | |
| CLM-1332-1.2 | Naphthalene (¹³ C ₆ , 99%) | *C ₆ C ₄ H ₈ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7425-1.2 | Naphthalene (unlabeled) | C ₁₀ H ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-3914-1.2 | DL-Nicotine (3',4',5'- ¹³ C ₃ , 99%) | CH ₃ *C ₃ CH ₂ NC ₅ H ₄ N | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9547-1.2 | Nicotine (unlabeled) | C ₁₀ H ₁₄ N ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-3913-S | 4-Nitrotoluene (ring- ¹³ C ₆ , 99%) | *C ₆ H ₄ CH ₃ NO ₂ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3891-1.2 | 4-Nitrotoluene (unlabeled) | C ₆ H ₄ CH ₃ NO ₂ | 1 mg/mL in acetonitrile | 1.2 mL |
| CLM-4811-1.2 | cis-Nonachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7445-1.2 | cis-Nonachlor (unlabeled) | C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4735-1.2 | trans-Nonachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7229-1.2 | trans-Nonachlor (unlabeled) | C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4306-1.2 | p-n-Nonylphenol (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ *C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-4306-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4559-1.2 | p-n-Nonylphenol (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-4559-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| CLM-4307-1.2 | p-n-Nonylphenol diethoxylate (ring- ¹³ C ₆ , 99%) | CH ₃ (CH ₂) ₈ *C ₆ H ₄ O(C ₂ H ₄ O) ₂ H | 100 µg/mL in nonane | 1.2 mL |
| CLM-4307-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |

(continued on next page)

Endocrine-Disrupting Compounds and Xenoestrogen Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|---|--|--|------------------|
| ULM-4521-1.2 | <i>p</i> -n-Nonylphenol diethoxylate (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ O(C ₂ H ₄ O) ₂ H | 100 µg/mL in nonane 100 µg/mL in methanol | 1.2 mL 1.2 mL |
| ULM-4521-M-1.2 | | | | |
| ULM-7147-1.2 | Nonylphenol diethoxylate-branched isomers (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ O(C ₂ H ₄ O) ₂ H | 100 µg/mL in nonane | 1.2 mL |
| CLM-4512-1.2 | <i>p</i> -n-Nonylphenol monoethoxylate | CH ₃ (CH ₂) ₈ *C ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-4512-M-1.2 | (ring- ¹³ C ₆ , 99%) | | 100 µg/mL in methanol | 1.2 mL |
| ULM-4520-1.2 | <i>p</i> -n-Nonylphenol monoethoxylate (unlabeled) | CH ₃ (CH ₂) ₈ C ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-4520-M-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-7146-1.2 | Nonylphenol monoethoxylate-branched isomers (unlabeled) | C ₉ H ₁₉ C ₆ H ₄ O(CH ₂) ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-4516-1.2 | <i>p</i> -n-Nonylphenol triethoxylate (ring- ¹³ C ₆ , 99%) CP 90% | CH ₃ (CH ₂) ₈ *C ₆ H ₄ O(C ₂ H ₄ O) ₃ H | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9679-1.2 | 1,1,1,3,10,12,12,12-Octachlorododecane (¹³ C ₁₂ , 99%) | *C ₁₂ H ₁₈ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9485-1.2 | 1,1,1,3,10,12,12,12-Octachlorododecane (unlabeled) | C ₁₂ H ₁₈ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4729-1.2 | Oxychlordane (¹³ C ₁₀ , 99%) | *C ₁₀ H ₄ Cl ₈ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-1.2 | Oxychlordane (unlabeled) | C ₁₀ H ₄ Cl ₈ O | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9849-1.2 | Benzyl paraben (benzyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₈ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9850-1.2 | Benzyl paraben (benzyl 4-hydroxybenzoate) (unlabeled) | C ₁₄ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| CLM-8285-1.2 | <i>n</i> -Butyl paraben (ring- ¹³ C ₆ , 99%) | *C ₆ C ₅ H ₁₄ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8287-1.2 | <i>n</i> -Butyl paraben (unlabeled) | C ₁₁ H ₁₄ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9761-1.2 | Ethyl paraben (ethyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₃ H ₁₀ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9760-1.2 | Ethyl paraben (ethyl 4-hydroxybenzoate) (unlabeled) | C ₉ H ₁₀ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9847-1.2 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₅ H ₁₄ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9848-1.2 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (unlabeled) | C ₁₁ H ₁₄ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9845-1.2 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9846-1.2 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (unlabeled) | C ₁₀ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| CLM-8249-1.2 | Methyl paraben (methyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₂ H ₈ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8250-1.2 | Methyl paraben (methyl 4-hydroxybenzoate) (unlabeled) | C ₈ H ₈ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9763-1.2 | <i>n</i> -Propyl paraben (<i>n</i> -propyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW ULM-9762-1.2 | <i>n</i> -Propyl paraben (<i>n</i> -propyl 4-hydroxybenzoate) (unlabeled) | C ₁₀ H ₁₂ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| DLM-2970-1.2 | Parathion (diethyl-D ₁₀ , 98%) | C ₁₀ D ₁₀ H ₄ NOPS | 100 µg/mL in nonane | 1.2 mL |
| ULM-8144-1.2 | Parathion (unlabeled) | C ₁₀ H ₁₄ NOPS | 100 µg/mL in nonane | 1.2 mL |
| CLM-7930-1.2 | Parlar 26 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7828-1.2 | Parlar 26 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8705-1.2 | Parlar 32 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8665-1.2 | Parlar 32 (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9005-1.2 | Parlar 38 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8719-1.2 | Parlar 39 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8767-1.2 | Parlar 39 (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9431-1.2 | Parlar 41 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9432-1.2 | Parlar 44 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7931-1.2 | Parlar 50 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7829-1.2 | Parlar 50 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7932-1.2 | Parlar 62 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7830-1.2 | Parlar 62 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |

Endocrine-Disrupting Compounds and Xenoestrogen Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--|--|---|---------------------------|------------|
| CLM-8720-1.2 | Parlar 69 ($^{13}\text{C}_{10}$, 99%) | $^*\text{C}_{10}\text{H}_9\text{Cl}_9$ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8768-1.2 | Parlar 69 (unlabeled) | $\text{C}_{10}\text{H}_9\text{Cl}_9$ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8721-1.2 | Parlar 70 ($^{13}\text{C}_{10}$, 99%) | $^*\text{C}_{10}\text{H}_9\text{Cl}_9$ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8769-1.2 | Parlar 70 (unlabeled) | $\text{C}_{10}\text{H}_9\text{Cl}_9$ | 10 µg/mL in nonane | 1.2 mL |
| EC-1404-3 | PCB-77 (3,3',4,4'-tetraCB) ($^{13}\text{C}_{12}$, 99%) | ($^*\text{C}_6\text{Cl}_2\text{H}_3$) ₂ | 40 µg/mL in nonane | 3 mL |
| EC-1425-3 | PCB-126 (3,3',4,4',5-pentaCB) ($^{13}\text{C}_{12}$, 99%) | $^*\text{C}_6\text{Cl}_3\text{H}_2 \cdot ^*\text{C}_6\text{Cl}_2\text{H}_3$ | 40 µg/mL in nonane | 3 mL |
| EC-1416-3 | PCB-169 (3,3',4,4',5,5'-hexaCB) ($^{13}\text{C}_{12}$, 99%) | ($^*\text{C}_6\text{Cl}_3\text{H}_2$) ₂ | 40 µg/mL in nonane | 3 mL |
| CLM-2050-1.2 | Pentachlorobenzene ($^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{HCl}_5$ | 100 µg/mL in isoctane | 1.2 mL |
| ULM-7234-1.2 | Pentachlorobenzene (unlabeled) | C_6HCl_5 | 100 µg/mL in isoctane | 1.2 mL |
| CLM-1955-1.2 | Pentachloronitrobenzene ($^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{Cl}_5\text{NO}_2$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7597-1.2 | Pentachloronitrobenzene (unlabeled) | $\text{C}_6\text{Cl}_5\text{NO}_2$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-661-1.2 | Pentachlorophenol ($^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{Cl}_5\text{OH}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6894-1.2 | Pentachlorophenol (unlabeled) | $\text{C}_6\text{Cl}_5\text{OH}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-8505-1.2 | Perfluoroctanesulfonate (PFOS), sodium salt ($^{13}\text{C}_8$, 99%) | $^*\text{C}_8\text{F}_{17}\text{NaO}_3\text{S}$ | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-9001-1.2 | Perfluoroctanesulfonate (PFOS), sodium salt (unlabeled) | $\text{C}_8\text{F}_{17}\text{NaO}_3\text{S}$ | 50 µg/mL in methanol | 1.2 mL |
| CLM-8005-1.2 | Perfluoroctanoic acid (PFOA) ($^{13}\text{C}_8$, 99%) | $^*\text{CF}_3(^*\text{CF}_2)_6^*\text{COOH}$ | 50 µg/mL in methanol | 1.2 mL |
| ULM-7451-1.2 | Perfluoroctanoic acid (PFOA) (unlabeled) (90:10 straight:branched isomers) CP 96% | $\text{CF}_3(\text{CF}_2)_6\text{COOH}$ | 50 µg/mL in methanol | 1.2 mL |
| CLM-7322-1.2 | cis-Permethrin (phenoxy- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_{15}\text{H}_{20}\text{Cl}_2\text{O}_3$ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8526-1.2 | cis-Permethrin (unlabeled) | $\text{C}_{21}\text{H}_{20}\text{Cl}_2\text{O}_3$ | 50 µg/mL in nonane | 1.2 mL |
| CLM-7323-1.2 | trans-Permethrin (phenoxy- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_{15}\text{H}_{20}\text{Cl}_2\text{O}_3$ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8527-1.2 | trans-Permethrin (unlabeled) | $\text{C}_{21}\text{H}_{20}\text{Cl}_2\text{O}_3$ | 50 µg/mL in nonane | 1.2 mL |
| CLM-2451-1.2 | Phenanthrene ($^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_8\text{H}_{10}$ | 100 µg/mL in nonane | 1.2 mL |
| DLM-371-1.2 | Phenanthrene (D_{10} , 98%) | $\text{C}_{14}\text{D}_{10}$ | 200 µg/mL in isoctane | 1.2 mL |
| ULM-7427-1.2 | Phenanthrene (unlabeled) | $\text{C}_{14}\text{H}_{10}$ | 200 µg/mL in isoctane | 1.2 mL |
| NEW DLM-695-1 | Phenol (ring- D_5 , 98%) | $\text{C}_6\text{D}_5\text{OH}$ | neat | 1 g |
| NEW DLM-7141-1.2 | Propoxur (isopropyl- D_7 , 98%) | $\text{C}_{11}\text{D}_7\text{H}_8\text{NO}_3$ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9765-1.2 | Propoxur (unlabeled) | $\text{C}_{11}\text{H}_{15}\text{NO}_3$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3739-1.2 | Simazine (ring- $^{13}\text{C}_3$, 99%) | $^*\text{C}_3\text{C}_4\text{H}_12\text{ClN}_5$ | 100 µg/mL in methanol | 1.2 mL |
| CLM-4694-1.2 | Tetrabromobisphenol A (ring- $^{13}\text{C}_{12}$, 99%) | ($^*\text{C}_6\text{Br}_2\text{H}_2\text{OH})_2\text{C}(\text{CH}_3)_2$ | 50 µg/mL in methanol | 1.2 mL |
| ULM-8734-1.2 | Tetrabromobisphenol A (unlabeled) | ($\text{C}_6\text{Br}_2\text{H}_2\text{OH})_2\text{C}(\text{CH}_3)_2$ | 50 µg/mL in methanol | 1.2 mL |
| NEW ULM-8734-T-1.2 | | | 50 µg/mL in toluene | 1.2 mL |
| ED-900 | 2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin ($^{13}\text{C}_{12}$, 99%) | ($^*\text{C}_6\text{H}_2\text{Cl}_2$) ₂ O ₂ | 50 µg/mL in nonane | 1.2 mL |
| ED-901 | 2,3,7,8-Tetrachlorodibenzo- <i>p</i> -dioxin (unlabeled) ($\text{C}_6\text{H}_2\text{Cl}_2$) ₂ O ₂ | ($\text{C}_6\text{H}_2\text{Cl}_2$) ₂ O ₂ | 50 µg/mL in nonane | 4 × 1.2 mL |
| DLM-7136-1.2 | Tributyltin chloride (D_{27} , 98%) | $\text{C}_{12}\text{D}_{27}\text{ClSn}$ | 100 µg/mL in MeCl | 1.2 mL |
| ULM-8061-1.2 | Tributyltin chloride (unlabeled) | $\text{C}_{12}\text{H}_{27}\text{ClSn}$ | 100 µg/mL in MeCl | 1.2 mL |
| CLM-4551-1.2 | 2,4,5-Trichlorophenoxyacetic acid (ring- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_2\text{H}_5\text{Cl}_3\text{O}_3$ | 100 µg/mL in MeCl | 1.2 mL |
| ULM-7213-1.2 | 2,4,5-Trichlorophenoxyacetic acid (unlabeled) | $\text{C}_8\text{H}_5\text{Cl}_3\text{O}_3$ | 100 µg/mL in MeCl | 1.2 mL |
| NEW CLM-9049-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (4,5,6- $^{13}\text{C}_3$, 99%) CP 97% | $^*\text{C}_5\text{C}_2\text{H}_2\text{Cl}_3\text{NO}$ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9204-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (unlabeled) | $\text{C}_5\text{H}_2\text{Cl}_3\text{NO}$ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-6779-1.2 | Tricosan (2',4,4'-Trichloro-2-hydroxydiphenyl ether) ($^{13}\text{C}_{12}$, 99%) | $^*\text{C}_{12}\text{H}_7\text{Cl}_3\text{O}_2$ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-6779-MT-1.2 | | | 100 µg/mL in MTBE | 1.2 mL |
| ULM-6935-1.2 | Tricosan (2',4,4'-Trichloro-2-hydroxydiphenyl ether) (unlabeled) | $\text{C}_{12}\text{H}_7\text{Cl}_3\text{O}_2$ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-6935-MT-1.2 | | | 100 µg/mL in MTBE | 1.2 mL |
| DLM-4479-1.2 | Trifluralin (di- <i>n</i> -propyl- D_{14} , 98%) | $\text{C}_{13}\text{D}_{14}\text{H}_2\text{F}_3\text{N}_3\text{O}_4$ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-4444-0.1 | Urethane (ethyl carbamate) (ethyl- D_5 , 98%) | $\text{C}_3\text{D}_5\text{H}_2\text{NO}_2$ | neat | 0.1 g |
| NEW DLM-167-C | Vinyl chloride (D_3 , 98%) | $\text{C}_2\text{D}_3\text{Cl}$ | 100 µg/mL in methanol-OD | 20 mL |
| NEW ULM-8224-1.2 | Vinyl chloride (unlabeled) | $\text{C}_2\text{H}_3\text{Cl}$ | 50 µg/mL in methanol | 1.2 mL |

This section represents only a partial listing of known and suspected endocrine-disrupting chemicals. If you do not see a standard listed here for a compound you are interested in, please contact CIL to discuss how we can help you with your research needs.

Industrial Chemical Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|---------------------------------------|--------|
| CLM-4674-1.2 | <i>n</i> -Butylbenzene (ring- ¹³ C ₆ , 99%) | *C ₆ H ₅ (CH ₂) ₃ CH ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4695-1.2 | 1,2-Dibromo-3-chloropropane (¹³ C ₃ , 99%) | *C ₃ H ₅ Br ₂ Cl | 100 µg/mL in methanol | 1.2 mL |
| CLM-6144-1.2 | 1,1-Dichloroethylene (random- ¹³ C, 99%) (stabilized with hydroquinone) | *CCH ₂ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-7214-1.2 | 1,1-Dichloroethylene (unlabeled) (stabilized with hydroquinone) | CCH ₂ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| CLM-6145-1.2 | 1,2-Dichloroethylene (¹³ C ₁ , 99%) (<i>cis/trans</i> mix) (stabilized with hydroquinone) | *CCH ₂ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-7215-1.2 | 1,2-Dichloroethylene (unlabeled) (<i>cis/trans</i> mix) (stabilized with hydroquinone) | CCH ₂ Cl ₂ | 100 µg/mL in methanol | 1.2 mL |
| CLM-1305-1.2 | 2,4-Dichlorophenol (¹³ C ₆ , 99%) | *C ₆ H ₃ Cl ₂ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-3374-1.2 | Epichlorohydrin (¹³ C ₃ , 99%) | *C ₃ H ₅ ClO | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-1008-1 | Epichlorohydrin (D ₅ , 98%) | CICD ₂ CD ₂ O | neat | 1 g |
| ULM-7403-1.2 | Epichlorohydrin (unlabeled) | CICH ₂ CHCH ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8008-1.2 | Hexachlorophene (¹³ C ₁₃ , 99%) | *CH ₂ [*C ₆ H(Cl) ₃ OH] ₂ | 50 µg/mL in methanol | 1.2 mL |
| ULM-8009-1.2 | Hexachlorophene (unlabeled) | CH ₂ [C ₆ H(Cl) ₃ OH] ₂ | 50 µg/mL in methanol | 1.2 mL |
| CLM-4745-1.2 | 4-Hydroxybenzoic acid (ring- ¹³ C ₆ , 99%) | *C ₆ CH ₆ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| ULM-8251-1.2 | 4-Hydroxybenzoic acid (unlabeled) | C ₇ H ₆ O ₃ | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-8792-1.2 | Sodium <i>bis</i> (2-ethylhexyl) sulfosuccinate (DOSS) (fumaric acid- ¹³ C ₄ , 99%) | *C ₄ C ₁₆ H ₃₇ NaO ₇ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-8807-1.2 | Sodium <i>bis</i> (2-ethylhexyl) sulfosuccinate (DOSS) (unlabeled) | C ₂₀ H ₃₇ NaO ₇ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8006-1.2 | Tetrachlorobisphenol A (ring- ¹³ C ₁₂ , 99%) | *C ₁₂ C ₃ H ₁₂ Cl ₄ O ₂ | 50 µg/mL in methanol | 1.2 mL |
| ULM-7606-1.2 | Tetrachlorobisphenol A (unlabeled) | C ₁₂ C ₃ H ₁₂ Cl ₄ O ₂ | 50 µg/mL in methanol | 1.2 mL |
| NEW DLM-9612-1.2 | Tetradecyl (tri- <i>n</i> -butyl) phosphonium bromide (D ₂₉ , 98%) | C ₂₆ H ₂₇ D ₂₉ PBr | 100 µg/mL in acetone:water (75:25) | 1.2 mL |
| NEW ULM-9609-1.2 | Tetradecyl (tri- <i>n</i> -butyl) phosphonium chloride (unlabeled) | C ₂₆ H ₅₆ ClP | 100 µg/mL in acetone:water (75:25) | 1.2 mL |
| DLM-7136-1.2 | Tributyltin chloride (D ₂₇ , 98%) | (C ₄ D ₉) ₃ ClSn | 100 µg/mL in MeCl-D ₂ | 1.2 mL |
| ULM-8061-1.2 | Tributyltin chloride (unlabeled) | (C ₄ H ₉) ₃ ClSn | 100 µg/mL in MeCl | 1.2 mL |
| CLM-6185-1.2 | 1,1,1-Trichloroethane (2- ¹³ C, 99%) | *CCH ₃ Cl ₃ | 100 µg/mL in methanol | 1.2 mL |
| DLM-2080-1.2 | 1,2,3-Trichloropropane (D ₅ , 98%) CP 95% | CD ₂ ClCDCICD ₂ Cl | 100 µg/mL in methanol | 1.2 mL |
| ULM-6911-1.2 | 1,2,3-Trichloropropane (unlabeled) | CH ₂ ClCHClCH ₂ Cl | 1 mg/mL in methanol | 1.2 mL |
| NEW CLM-9095 | Trimethylolpropane phosphate (3,4,9,10- ¹³ C ₄ , 99%) CP 95% | *C ₄ C ₂ H ₁₁ O ₄ P | Inquire | |
| NEW ULM-9096 | Trimethylolpropane phosphate (unlabeled) CP 95% | C ₆ H ₁₁ O ₄ P | Inquire | |
| NEW DLM-167-C | Vinyl chloride (D ₃ , 98%) | C ₂ D ₃ Cl | 100 µg/mL in methanol-OD | 20 mL |
| NEW ULM-8224-1.2 | Vinyl chloride (unlabeled) | C ₂ H ₃ Cl | 50 µg/mL in methanol | 1.2 mL |

Explosives Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------------|--|--|---|---------------|
| CLM-1519-S | 1,3-Dinitrobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile neat | 1 mL 0.1 g |
| CLM-1519-0.1 | | | | |
| ULM-3850-1.2 | 1,3-Dinitrobenzene (unlabeled) | $\text{C}_6\text{H}_4(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile | 1.2 mL |
| DLM-299-10 | 2,4-Dinitrophenol (ring-D ₃ , 98%) (contains 0.35 mg/mL deuterium oxide) | (NO_2) ₂ $\text{C}_6\text{D}_3\text{OH}$ | 1 mg/mL in methanol-OD | 10 mL |
| NEW ULM-8706-10 | 2,4-Dinitrophenol (unlabeled) (contains 0.35 mg/mL water) | (NO_2) ₂ $\text{C}_6\text{H}_3\text{OH}$ | 1 mg/mL in methanol | 10 mL |
| DLM-2207-S | 2,4-Dinitrotoluene (ring-D ₃ , 98%) | $\text{C}_6\text{D}_3\text{CH}_3(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3888-S | 2,4-Dinitrotoluene (unlabeled) | $\text{C}_6\text{H}_3\text{CH}_3(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile | 1 mL |
| DLM-1939-S | 2,6-Dinitrotoluene (methyl-D ₃ , 98%) | $\text{C}_6\text{H}_3\text{CD}_3(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3889-S | 2,6-Dinitrotoluene (unlabeled) | $\text{C}_6\text{H}_3\text{CH}_3(\text{NO}_2)_2$ | 1 mg/mL in acetonitrile | 1 mL |
| CNLM-7963-S | HMX ($^{13}\text{C}_4$, 99%; ring- $^{15}\text{N}_4$, 98%) | * $\text{C}_4\text{H}_8\text{N}_4^*\text{N}_4\text{O}_8$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-7969-1 | HMX (unlabeled) | $\text{C}_4\text{H}_8\text{N}_4\text{N}_4\text{O}_8$ | 1 mg/mL in acetonitrile | 1 mL |
| NEW CLM-675 | Nitrobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_5\text{NO}_2$ | Inquire | |
| NEW DLM-294-5 | Nitrobenzene (D ₅ , 99%) | $\text{C}_6\text{D}_5\text{NO}_2$ | neat | 5 g |
| NEW DLM-294-10 | | | neat | 10 g |
| ULM-3892-1.2 | Nitrobenzene (unlabeled) | $\text{C}_6\text{H}_5\text{NO}_2$ | 1 mg/mL in acetonitrile | 1.2 mL |
| NEW NLM-814-1.2 | Nitroglycerin (trinitroglycerol) ($^{15}\text{N}_3$, 98%) | $\text{C}_3\text{H}_5(*\text{NO}_3)_3$ | 1 mg/mL in ethanol | 1.2 mL |
| ULM-3893-S | Nitroglycerin (trinitroglycerol) (unlabeled) | $\text{C}_3\text{H}_5(\text{NO}_3)_3$ | 1 mg/mL in acetonitrile | 1 mL |
| CLM-3912-S | 2-Nitrotoluene (ring- $^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{CH}_3\text{NO}_2$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3890-1.2 | 2-Nitrotoluene (unlabeled) | $\text{C}_6\text{H}_4\text{CH}_3\text{NO}_2$ | 1 mg/mL in acetonitrile | 1.2 mL |
| CLM-3913-S | 4-Nitrotoluene (ring- $^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{CH}_3\text{NO}_2$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3891-1.2 | 4-Nitrotoluene (unlabeled) | $\text{C}_6\text{H}_4\text{CH}_3\text{NO}_2$ | 1 mg/mL in acetonitrile | 1.2 mL |
| CNLM-7987-S | RDX ($^{13}\text{C}_3$, 99%; $^{15}\text{N}_3$, 98%) | * $\text{C}_3\text{H}_6\text{N}_3(*\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1 mL |
| CLM-3846-S | RDX ($^{13}\text{C}_3$, 99%) | * $\text{C}_3\text{H}_6\text{N}_3(\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3847-S | RDX (unlabeled) | $\text{C}_3\text{H}_6\text{N}_3(\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1.2 mL |
| CLM-3848-S | 1,3,5-Trinitrobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_3(\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1 mL |
| ULM-3849-1.2 | 1,3,5-Trinitrobenzene (unlabeled) | $\text{C}_6\text{H}_3(\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1.2 mL |
| CNLM-3643-S | 2,4,6-Trinitrotoluene (TNT) ($^{13}\text{C}_7$, 99%; $^{15}\text{N}_3$, 98%) | * $\text{C}_7\text{H}_5(*\text{NO}_2)_3$ | 1 mg/mL in benzene (wetted with H ₂ O 33% by weight) | 1 mL |
| ULM-3845-1.2 | 2,4,6-Trinitrotoluene (TNT) (unlabeled) | $\text{C}_7\text{H}_5(\text{NO}_2)_3$ | 1 mg/mL in acetonitrile | 1.2 mL |

Note: Shipping restrictions on explosive compounds may prevent CIL from shipping certain standards, especially outside of the US. Please contact CIL to confirm availability of these explosive standards.

***n*-Alkane Standards**

| Catalog No. | Compound | Formula | Amount |
|-------------------------|--|--|--------|
| DLM-1213-1 | <i>n</i> -Pentane (D ₁₂ , 98%) | CD ₃ (CD ₂) ₃ CD ₃ | 1 g |
| DLM-1213-5 | | | 5 g |
| DLM-139-1 | <i>n</i> -Hexane (D ₁₄ , 98%) | CD ₃ (CD ₂) ₄ CD ₃ | 1 g |
| DLM-139-5 | | | 5 g |
| DLM-423-1 | <i>n</i> -Heptane (D ₁₆ , 98%) | CD ₃ (CD ₂) ₅ CD ₃ | 1 g |
| DLM-423-5 | | | 5 g |
| DLM-50-1 | <i>n</i> -Octane (D ₁₈ , 99%) | CD ₃ (CD ₂) ₆ CD ₃ | 1 g |
| DLM-50-5 | | | 5 g |
| DLM-2438-1 | <i>n</i> -Nonane (D ₂₀ , 98%) | CD ₃ (CD ₂) ₇ CD ₃ | 1 g |
| DLM-2438-5 | | | 5 g |
| DLM-133-1 | <i>n</i> -Decane (D ₂₂ , 99%) | CD ₃ (CD ₂) ₈ CD ₃ | 1 g |
| DLM-133-5 | | | 5 g |
| DLM-338-1 | <i>n</i> -Dodecane (D ₂₆ , 98%) | CD ₃ (CD ₂) ₁₀ CD ₃ | 1 g |
| DLM-338-5 | | | 5 g |
| NEW DLM-1354-0.1 | <i>n</i> -Tridecane (D ₂₈ , 98%) | CD ₃ (CD ₂) ₁₁ CD ₃ | 0.1 g |
| DLM-1354-0.5 | | | 0.5 g |
| DLM-670-1 | <i>n</i> -Tetradecane (D ₃₀ , 98%) | CD ₃ (CD ₂) ₁₂ CD ₃ | 1 g |
| DLM-670-5 | | | 5 g |
| DLM-1283-1 | <i>n</i> -Pentadecane (D ₃₂ , 98%) | CD ₃ (CD ₂) ₁₃ CD ₃ | 1 g |
| DLM-1283-5 | | | 5 g |
| DLM-203-0.1 | <i>n</i> -Hexadecane (D ₃₄ , 98%) | CD ₃ (CD ₂) ₁₄ CD ₃ | 0.1 g |
| NEW DLM-203-1 | | | 1 g |
| DLM-203-5 | | | 5 g |
| NEW DLM-1342-1 | <i>n</i> -Heptadecane (D ₃₆ , 98%) CP 95% | CD ₃ (CD ₂) ₁₅ CD ₃ | 1 g |
| DLM-1342-5 | | | 5 g |
| DLM-1346-0.1 | <i>n</i> -Nonadecane (D ₄₀ , 98%) | CD ₃ (CD ₂) ₁₇ CD ₃ | 0.1 g |
| DLM-1346-1 | | | 1 g |
| DLM-2208-0.5 | <i>n</i> -Eicosane (D ₄₂ , 98%) | CD ₃ (CD ₂) ₁₈ CD ₃ | 0.5 g |
| DLM-2208-1 | | | 1 g |
| DLM-3336-1 | <i>n</i> -Tricosane (D ₄₈ , 98%) | CD ₃ (CD ₂) ₂₁ CD ₃ | 1 g |
| DLM-2209-0.5 | <i>n</i> -Tetracosane (D ₅₀ , 98%) | CD ₃ (CD ₂) ₂₂ CD ₃ | 0.5 g |
| DLM-2210-0.5 | <i>n</i> -Triacontane (D ₆₂ , 98%) | CD ₃ (CD ₂) ₂₈ CD ₃ | 0.5 g |
| DLM-2724-1 | <i>n</i> -Dotriacontane (D ₆₆ , 98%) | CD ₃ (CD ₂) ₃₀ CD ₃ | 1 g |
| DLM-2634-1 | <i>n</i> -Hexatriacontane (D ₇₄ , 98%) | CD ₃ (CD ₂) ₃₄ CD ₃ | 1 g |

Priority Pollutant Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------|---|-------------------------------|--------------------------------|--------|
| DLM-9-10 | Acetone (D_6 , 99.9%) | CD_3COCD_3 | neat | 10 g |
| CLM-856-0.1 | Acrylonitrile ($^{13}C_3$, 99%) (inhibited with 0.1% 4-methoxy phenol) | $H_2^*C=^*CH^*CN$ | neat | 0.1 g |
| DLM-820-1 | Acrylonitrile (D_3 , 98%) | $D_2C=CDCN$ | neat | 1 g |
| DLM-820-5 | (inhibited with 0.1% 4-methoxy phenol) | | neat | 5 g |
| DLM-2030-1.2 | 2-Aminonaphthalene (ring- D_7 , 98%) | $C_{10}D_7NH_2$ | 1 mg/mL in benzene | 1.2 mL |
| ULM-9376-1.2 | 2-Aminonaphthalene (unlabeled) | $C_{10}H_7NH_2$ | 1 mg/mL in benzene | 1.2 mL |
| DLM-7658 | 1-Amino-2-propanol (D_9 , 98%) | C_3D_9NO | Inquire | |
| CLM-714-0.1 | Aniline ($^{13}C_6$, 99%) | $*C_6H_5NH_2$ | neat | 0.1 g |
| CLM-714-0.25 | | | neat | 0.25 g |
| DLM-862-1 | Aniline (ring- D_5 , 98%) | $C_6D_5NH_2$ | neat | 1 g |
| DLM-862-5 | | | neat | 5 g |
| DLM-106-5 | Aniline (D_7 , 98%) | $C_6D_5ND_2$ | neat | 5 g |
| CLM-182-0.1 | Benzene ($^{13}C_6$, 99%) | $*C_6H_6$ | neat | 0.1 g |
| CLM-182-0.5 | | | neat | 0.5 g |
| DLM-1-5 | Benzene (D_6 , 99.5%) | C_6D_6 | neat | 5 g |
| CDLM-629-0.1 | Benzene ($^{13}C_6$, 99%; D_6 , 98%) | $*C_6D_6$ | neat | 0.1 g |
| DLM-1338-1.2 | Benzidine (ring- D_8 , 98%) | $C_{12}D_8(NH_2)_2$ | 100 μ g/mL in toluene | 1.2 mL |
| DLM-122-1 | Benzoic acid (ring- D_5 , 98%) | $C_8D_5CO_2H$ | neat | 1 g |
| DLM-122-5 | | | neat | 5 g |
| DLM-1663-1 | 1,4-Benzoquinone (D_4 , 98%) | $O(C_6D_4)O$ | neat | 1 g |
| CLM-3235-1.2 | Biphenyl ($^{13}C_{12}$, 99%) | $*C_{12}H_{10}$ | 100 μ g/mL in nonane | 1.2 mL |
| DLM-494-1 | Biphenyl (D_{10} , 98%) | $C_{12}D_{10}$ | neat | 1 g |
| DLM-494-5 | | | neat | 5 g |
| ULM-1710-1.2 | Biphenyl (unlabeled) | $C_{12}H_{10}$ | 50 μ g/mL in nonane | 1.2 mL |
| ULM-1710-0.5 | | | neat | 0.5 g |
| DLM-1945-0.1 | Bis(2-chloroethoxy) methane (chloroethoxy- D_8 , 98%) | $CH_2(OCD_2CD_2Cl)_2$ | neat | 0.1 g |
| DLM-2004-0.05 | Bis(2-chloroethyl) ether (D_8 , 98%) | $ClCD_2CD_2OCD_2CD_2Cl$ | neat | 0.05 g |
| DLM-2004-0.1 | | | neat | 0.1 g |
| DLM-2138 | Bis(2-chloroisopropyl) ether (D_{12} , 95%) | $C_6D_{12}C_{12}O$ | Inquire | |
| ULM-3693 | Bis(2-chloroisopropyl) ether (unlabeled) | $C_6H_{12}C_{12}O$ | Inquire | |
| CLM-4325-1.2 | Bisphenol A (ring- $^{13}C_{12}$, 99%) | $*(C_6H_4OH)_2C(CH_3)_2$ | 100 μ g/mL in acetonitrile | 1.2 mL |
| ULM-7106-1.2 | Bisphenol A (unlabeled) | $(C_6H_4OH)_2C(CH_3)_2$ | 100 μ g/mL in acetonitrile | 1.2 mL |
| ULM-8654-1.2 | 2,4'-Bisphenol A (2-(2-hydroxyphenyl)-2-(4-hydroxyphenyl)propane) (unlabeled) | $(C_6H_4OH)_2C(CH_3)_2$ | 100 μ g/mL in acetonitrile | 1.2 mL |
| DLM-872-0.1 | Bromochloromethane (D_2 , 98%) | CD_2ClBr | neat | 0.1 g |
| CLM-2090-1 | Bromodichloromethane (^{13}C , 99%) (stabilized with K_2CO_3) | Br^*CHCl_2 | neat | 1 g |
| ULM-8480 | Bromodichloromethane (unlabeled) | $BrCHCl_2$ | Inquire | |
| DLM-874-10 | Bromoethane (D_5 , 98%) | CD_3CD_2Br | neat | 10 g |
| DLM-103-1 | 2-Bromoethanol (1,1,2,2- D_4 , 98%) CP 95%+ | $BrCD_2CD_2OH$ | neat | 1 g |
| DLM-103-5 | | | neat | 5 g |
| CLM-726-0.1 | Bromoform (^{13}C , 99%) (stabilized with copper wire) | $*CHBr_3$ | neat | 0.1 g |
| CLM-726-0.5 | | | neat | 0.5 g |
| DLM-400-10 | Bromoform (D , 99.5%) (stabilized with copper wire) | $CDBr_3$ | neat | 10 g |
| DLM-400-25 | | | neat | 25 g |
| CLM-1217-1 | Bromomethane (^{13}C , 99%) * | $*CH_3Br$ | neat | 1 L |
| DLM-401-5 | Bromomethane (D_3 , 99%) * | CD_3Br | neat | 5 g |
| DLM-1910-0.1 | 2-Butanone (4,4,4- D_3 , 98%) | $CD_3CH_2COCH_3$ | neat | 0.1 g |
| DLM-1910-1 | | | neat | 1 g |
| DLM-663-0.1 | 2-Butanone (1,1,1,3,3- D_5 , 98%) | $CH_3CD_2COCD_3$ | neat | 0.1 g |
| DLM-663-1 | | | neat | 1 g |
| DLM-663-5 | | | neat | 5 g |
| NEW DLM-8811-1.2 | 2-Butoxyethanol (1,1,2,2,- D_4 , 99%) | $CH_3CH_2CH_2CH_2OCD_2CD_2OH$ | 1000 μ g/mL in water | 1.2 mL |
| NEW ULM-9046-1.2 | 2-Butoxyethanol (unlabeled) | $C_6H_{14}O_2$ | 1000 μ g/mL in water | 1.2 mL |
| DLM-2134-0.1 | Carbazole (ring- D_8 , 98%) | $C_{12}D_8NH$ | neat | 0.1 g |

*Gases require a breakseal flask or cylinder and valve at an additional charge. Breakseal flasks are only available for certain gases at atmospheric pressure.

(continued on next page)

Priority Pollutant Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------------|--|---|---------------------------------|---------|
| CLM-731-0.1 | Carbon tetrachloride (^{13}C , 99%) | * CCl_4 | neat | 0.1 g |
| CLM-731-0.5 | | | neat | 0.5 g |
| CLM-731-1 | | | neat | 1 g |
| CLM-1520-1 | Catechol ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4(\text{OH})_2$ | neat | 1 mg |
| DLM-1912-5 | Catechol (D_6 , 98%) | $\text{C}_6\text{D}_4(\text{OD})_2$ | neat | 5 g |
| CLM-2284-1 | 4-Chlorocatechol ($^{13}\text{C}_6$, 99%) | $\text{Cl}^*\text{C}_6\text{H}_3(\text{OH})_2$ | neat | 1 mg |
| ULM-1701-0.1 | 4-Chlorocatechol (unlabeled) CP 90-95% | $\text{CIC}_6\text{H}_3(\text{OH})_2$ | neat | 0.1 g |
| CLM-2091 | Chlorodibromomethane (^{13}C , 99%) | Br_2^*CHCl | | Inquire |
| NEW DLM-1171-A-1.2 | Chloroethane (D_5 , 98%) | $\text{CD}_3\text{CD}_2\text{Cl}$ | 1000 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| | | | neat | 5 g |
| DLM-1928-0.5 | 2-Chloroethanol (1,1,2,2-D ₄ , 98%) | $\text{CICD}_2\text{CD}_2\text{OH}$ | neat | 0.5 g |
| CLM-262-0.1 | Chloroform (^{13}C , 99%) | * CHCl_3 | neat | 0.1 g |
| CLM-262-0.5 | | | neat | 0.05 g |
| CLM-262-1 | | | neat | 1 g |
| ULM-1705-0.1 | 4-Chloroguaiaacol (unlabeled) CP 85-90% | $\text{CIC}_6\text{H}_3(\text{OH})(\text{OCH}_3)$ | neat | 0.1 g |
| DLM-2037-1 | Chloroiodomethane (D_2 , 98%) (stabilized with copper wire) | CICD_2I | neat | 1 g |
| CLM-339-1 | Chloromethane (^{13}C , 99%) | * CH_3Cl | neat | 1 L |
| DLM-337-1-BS | Chloromethane (D_3 , 99%) | CD_3Cl | neat | 1 L |
| DLM-337-1-LB | Chloromethane (D_3 , 99%) | CD_3Cl | neat | 1 L |
| DLM-2205-0.01 | 4-Chloro-3-methylphenol (ring-2,6-D ₂ , 98%) | $\text{C}_7\text{D}_2\text{H}_4\text{ClO}$ | neat | 0.01 g |
| DLM-2205-0.1 | | | neat | 0.1 g |
| DLM-2005-1.2 | 2-Chloronaphthalene (D_7 , 98%) | $\text{C}_{10}\text{D}_7\text{Cl}$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| DLM-2005-0.01 | | | neat | 0.01 g |
| DLM-2005-0.1 | | | neat | 0.1 g |
| CLM-1559-1 | 4-Chloronitrobenzene ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_4\text{NO}_2\text{Cl}$ | neat | 1 mg |
| DLM-1930-0.1 | 4-Chlorophenyl phenyl ether (phenyl-D ₅ , 98%) | $\text{CIC}_6\text{H}_4\text{OC}_6\text{D}_5$ | neat | 0.1 g |
| ULM-2421-0.1 | 4-Chlorophenyl phenyl ether (unlabeled) | $\text{CIC}_6\text{H}_4\text{OC}_6\text{H}_5$ | neat | 0.1 g |
| DLM-3014-1 | 2-Chloropropene (D_5 , 98%) | $\text{D}_3\text{CClC=CD}_2$ | neat | 1 g |
| DLM-3014-5 | | | neat | 5 g |
| NEW DLM-3016-1 | <i>o</i> -Cresol (D_8 , 98%) | $\text{D}_3\text{CC}_6\text{D}_4\text{OD}$ | neat | 1 g |
| | | | neat | 5 g |
| CLM-7341 | <i>p</i> -Cresol (ring- $^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{CH}_8\text{O}$ | | Inquire |
| NEW DLM-3017-1 | <i>p</i> -Cresol (D_8 , 98%) | $\text{D}_3\text{CC}_6\text{D}_4\text{OD}$ | neat | 1 g |
| | | | neat | 5 g |
| DLM-1386-1 | Decalin (D_{18} , 99%) (<i>cis/trans</i> mixture) | $\text{C}_{10}\text{D}_{18}$ | neat | 1 g |
| DLM-1386-5 | | | neat | 5 g |
| DLM-1843-5 | <i>trans</i> -Decalin (D_{18} , 98%) | $\text{C}_{10}\text{D}_{18}$ | neat | 5 g |
| CLM-1544-1.2 | Dibenzo- <i>p</i> -dioxin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{O}_2$ | 50 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-1711-1.2 | Dibenzo- <i>p</i> -dioxin (unlabeled) | $\text{C}_{12}\text{H}_8\text{O}_2$ | 50 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-1711-0.01 | | | neat | 0.01 g |
| CLM-1561-1.2 | Dibenzofuran ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{O}$ | 50 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| DLM-2276-0.05 | Dibenzofuran (D_8 , 98%) | $\text{C}_{12}\text{D}_8\text{O}$ | neat | 0.05 g |
| ULM-1712-1.2 | Dibenzofuran (unlabeled) | $\text{C}_{12}\text{H}_8\text{O}$ | 50 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-1712-0.05 | | | neat | 0.05 g |
| DLM-2206-0.1 | Dibenzothiophene (D_8 , 98%) | $\text{C}_{12}\text{D}_8\text{S}$ | neat | 0.1 g |
| CLM-483-0.1 | 1,2-Dibromoethane ($^{13}\text{C}_2$, 99%) | $\text{Br}^*\text{CH}_2\text{CH}_2\text{Br}$ | neat | 0.1 g |
| CLM-483-1 | | | neat | 1 g |
| CLM-735-1 | 3,4-Dichloroaniline ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_3\text{Cl}_2\text{NH}_2$ | neat | 1 mg |
| DLM-3022-1.2 | 3,3'-Dichlorobenzidine (ring-D ₆ , 98%) | $\text{C}_{12}\text{D}_6\text{N}_2\text{Cl}_2$ | 1 mg/mL in benzene | 1.2 mL |
| ULM-1702-0.1 | 4,5-Dichlorocatechol (unlabeled) CP 95-99% | $\text{Cl}_2\text{C}_6\text{H}_3(\text{OH})_2$ | neat | 0.1 g |
| DLM-1934-0.1 | 1,1-Dichloroethane (2,2,2-D ₃ , 98%) | CD_3CHCl_2 | neat | 0.1 g |
| DLM-1934-0.25 | | | neat | 0.25 g |
| DLM-18-1 | 1,2-Dichloroethane (D_4 , 99%) | $\text{CICD}_2\text{CD}_2\text{Cl}$ | neat | 1 g |
| DLM-18-5 | | | neat | 5 g |
| DLM-1935-0.1 | 1,1-Dichloroethylene (2,2-D ₂ , 98%) (inhibited with hydroquinone) | $\text{CD}_2=\text{CCl}_2$ | neat | 0.1 g |
| DLM-1935-1 | | | neat | 1 g |

*Gases require a breakseal flask or cylinder and valve at an additional charge. Breakseal flasks are only available for certain gases at atmospheric pressure.

Priority Pollutant Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|---|---|---------------------------|--------|
| DLM-1936-0.1 | 1,2-Dichloroethylene (1,2-D ₂ , 98%) (<i>cis/trans</i> mixture) | CICD=CDCl | neat | 0.1 g |
| DLM-1936-1 | | | neat | 1 g |
| DLM-1937-0.1 | 1,2-Dichloropropane (D ₆ , 98%) | CICD ₂ CDClCD ₃ | neat | 0.1 g |
| DLM-1937-0.25 | | | neat | 0.25 g |
| DLM-2112-1.2 | 1,3-Dichloro-2-propanol (D ₅ , 98%) | CICD ₂ CD(OH)CD ₂ Cl | 1 mg/mL in methanol | 1.2 mL |
| ULM-8092-1.2 | 1,3-Dichloro-2-propanol (unlabeled) | CICH ₂ CH(OH)CH ₂ Cl | 1 mg/mL in methanol | 1.2 mL |
| DLM-1938-0.1 | 1,3-Dichloropropene (D ₄ , 98%) (<i>cis/trans</i> mixture) | CICD ₂ CD=CDCl | neat | 0.1 g |
| ULM-1700-0.1 | 5,6-Dichlorovanillin (unlabeled) | Cl ₂ C ₆ H(CHO)(OH)(OCH ₃) | neat | 0.1 g |
| DLM-1592-1 | Diethyl ether (D ₁₀ , 99%) | O(CD ₃ CD ₂) ₂ | neat | 1 g |
| DLM-1592-5 | | | neat | 5 g |
| ULM-8235-1.2 | Diethylene glycol (unlabeled) | O(CH ₂ CH ₂ OH) ₂ | 1 mg/mL in methanol | 1.2 mL |
| CLM-1006-0.5 | Diodomethane (¹³ C, 99%) (stabilized with copper wire) | *CH ₂ I ₂ | neat | 0.5 g |
| DLM-3190-1 | N,N-Dimethylaniline (D ₁₁ , 98%) | C ₆ D ₅ N(CD ₃) ₂ | neat | 1 g |
| CLM-503-0.5 | N,N-Dimethylformamide (carbonyl- ¹³ C, 99%) | H*CON(CH ₃) ₂ | neat | 0.5 g |
| CLM-503-1 | | | neat | 1 g |
| DLM-1366-1.2 | Dimethyl phthalate (ring-D ₄ , 98%) | C ₆ D ₄ -1,2-(CO ₂ CH ₃) ₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1366-0.1 | | | neat | 0.1 g |
| DLM-3024-5 | 1,3-Dinitrobenzene (D ₄ , 98%) | C ₆ D ₄ N ₂ O ₄ | neat | 5 g |
| DLM-3173-0.1 | 4,6-Dinitro-2-methylphenol (ring-D ₂ , 98%) | CH ₃ C ₆ D ₂ (NO ₂) ₂ OH | neat | 0.1 g |
| DLM-299-10 | 2,4-Dinitrophenol (ring-D ₃ , 98%) | C ₆ D ₃ (NO ₂) ₂ OH | 1 mg/mL in methanol-OD | 10 mL |
| NEW ULM-8706-10 | 2,4-Dinitrophenol (unlabeled) (contains 0.35 mg/mL water) | (NO ₂) ₂ C ₆ H ₃ OH | 1 mg/mL in methanol | 10 mL |
| DLM-2207-S | 2,4-Dinitrotoluene (ring-D ₃ , 98%) | H ₃ CC ₆ D ₃ (NO ₂) ₂ | 1 mg/mL in acetonitrile | 1 mL |
| DLM-1939-S | 2,6-Dinitrotoluene (methyl-D ₃ , 98%) | D ₃ CC ₆ H ₅ (NO ₂) ₂ | 1 mg/mL in acetonitrile | 1 mL |
| DLM-28-SM-1.2 | 1,4-Dioxane (<i>p</i> -dioxane) (D ₈ , 99%) | C ₄ D ₈ O ₂ | 1 mg/mL in methanol | 1.2 mL |
| DLM-28-5 | | | neat | 5 g |
| DLM-28-10 | | | neat | 10 g |
| DLM-28-25 | | | neat | 25 g |
| ULM-7840-1.2 | 1,4-Dioxane (<i>p</i> -dioxane) (unlabeled) | C ₄ H ₈ O ₂ | 1 mg/mL in methanol | 1.2 mL |
| DLM-2133-0.1 | Diphenylamine (diphenyl-D ₁₀ , 98%) | C ₆ D ₅ NHC ₆ D ₅ | neat | 0.1 g |
| CLM-1587-1.2 | Diphenyl ether (¹³ C ₁₂ , 99%) | (*C ₆ H ₅) ₂ O | 50 µg/mL in nonane | 1.2 mL |
| DLM-2211-0.1 | Diphenyl ether (D ₁₀ , 98%) | (C ₆ D ₅) ₂ O | neat | 0.1 g |
| DLM-3026-0.05 | 1,2-Diphenylhydrazine (diphenyl-D ₁₀ , 98%) | C ₁₂ D ₁₀ H ₅ N ₂ | neat | 0.05 g |
| DLM-3026-0.1 | | | neat | 0.1 g |
| NEW DLM-4880-1.2 | N,N'-Diphenyl- <i>p</i> -phenylenediamine (D ₁₄ , 98%) CP 95% | C ₆ D ₅ NHC ₆ D ₄ NHC ₆ D ₅ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9465-1.2 | N,N'-Diphenyl- <i>p</i> -phenylenediamine (unlabeled) | C ₆ H ₅ NHC ₆ H ₄ NHC ₆ H ₅ | 100 µg/mL in nonane | 1.2 mL |
| DLM-411-5 | Durene (1,2,4,5-tetramethylbenzene) (D ₁₄ , 98%) | C ₆ D ₂ (CD ₃) ₄ | neat | 5 g |
| CLM-3374-1.2 | Epichlorohydrin (¹³ C ₃ , 99%) | *C ₃ H ₅ ClO | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-1008-1 | Epichlorohydrin (D ₅ , 98%) | CICD ₂ CDClD ₂ O | neat | 1 g |
| ULM-7403-1.2 | Epichlorohydrin (unlabeled) | CICH ₂ CHCH ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-686-5 | Ethylbenzene (ethyl-D ₅ , 98%) | C ₆ H ₅ CD ₂ CD ₃ | neat | 5 g |
| DLM-199-10 | Ethylbenzene (D ₁₀ , 98%) | C ₆ D ₅ CD ₂ CD ₃ | neat | 10 g |
| DLM-4304-10 | Ethylbenzene (D ₁₀ , 99%) | C ₆ D ₅ CD ₂ CD ₃ | neat | 10 g |
| CLM-473-0.1 | Ethylene oxide (¹³ C ₂ , 99%) * (airfreight forbidden) | *CH ₂ *CH ₂ O | neat | 0.1 g |
| CLM-473-0.5 | | | neat | 0.5 g |
| DLM-271-5 | Ethylene oxide (D ₄ , 99%) * (airfreight forbidden) | CD ₂ CD ₂ O | neat | 5 g |
| CLM-810-1 | Guaiacol (ring- ¹³ C ₆ , 99%) | CH ₃ O*C ₆ H ₄ OH | neat | 1 mg |
| CLM-2145-1.2 | Hexachloro-1,3-butadiene (¹³ C ₄ , 99%) | *CCl ₂ =*CCl*CCl=*CCl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-2145-0.01 | | | neat | 0.01 g |
| NEW ULM-7526-1.2 | Hexachloro-1,3-butadiene (unlabeled) | CCl ₂ =CClCCl=CCl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| CLM-2110-5 | Hexachlorocyclopentadiene (¹³ C ₄ , 99%) | *C ₄ CCl ₆ | neat | 5 mg |
| CLM-2110-10 | | | neat | 10 mg |
| CLM-2003-0.1 | Hexachloroethane (1- ¹³ C, 99%) | CCl ₃ *CCl ₃ | neat | 0.1 g |
| CLM-2003-0.5 | | | neat | 0.5 g |
| ULM-6074-60 | 1,2,4,5,7,8-Hexachloroxanthene (unlabeled) | C ₁₃ H ₄ Cl ₆ O | neat | 60 µg |
| DLM-277-0.1 | Hexanoic acid (D ₁₁ , 98%) | CD ₃ (CD ₂) ₄ CO ₂ H | neat | 0.1 g |
| DLM-277-1 | | | neat | 1 g |

(continued on next page)

Priority Pollutant Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|---|---------------------------|-----------|
| DLM-1522-1 | Hydroquinone (ring-D ₄ , 98%) | HOC ₆ D ₄ OH | neat | 1 g |
| NEW NLM-6715-1.2 | 8-Hydroxy-2'-deoxyguanosine (¹⁵ N ₅ , 98%) CP 95% | C ₁₀ H ₁₃ *N ₅ O ₅ | 25 µg/mL in water | 1.2 mL |
| NEW ULM-9700-1.2 | 8-Hydroxy-2'-deoxyguanosine (unlabeled) | C ₁₀ H ₁₃ N ₅ O ₅ | 25 µg/mL in water | 1.2 mL |
| NEW CLM-9593 | 3-Hydroxymethyltetrahydrofuran (¹³ C ₅ , 99%) | *C ₅ H ₁₀ O ₂ | | Inquire |
| ULM-2-4X25 | Isooctane (unlabeled) | (CH ₃) ₃ CCH ₂ CH(CH ₃) ₂ | neat | 4 × 25 mL |
| DLM-1943-0.1 | Isophorone (3-methyl-D ₃ ; 2,4,4,6,6-D ₅ , 98%) | CD ₂ C(CH ₃) ₂ CD ₂ C(CD ₃)=CDCO | neat | 0.1 g |
| CLM-7864-1.2 | Leucomalachite green (phenyl- ¹³ C ₆ , 99%) | *C ₆ H ₅ CH[C ₆ H ₄ N(CH ₃) ₂] ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7870-1.2 | Leucomalachite green (unlabeled) | C ₆ H ₅ CH[C ₆ H ₄ N(CH ₃) ₂] ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-24-5 | Methanol (D ₄ , 99.8%) | CD ₃ OD | neat | 5 g |
| DLM-24-10 | | | neat | 10 g |
| CLM-1593-0.25 | Methylene chloride (¹³ C, 99%) | *CH ₂ Cl ₂ | neat | 0.25 g |
| CLM-1593-0.5 | | | neat | 0.5 g |
| DLM-23-5 | Methylene chloride (D ₂ , 99.9%) | CD ₂ Cl ₂ | neat | 5 g |
| DLM-2277-1 | 2-(4-Methylphenyl) propane (D ₁₄ , 98%) | D ₃ CC ₆ D ₄ CD(CD ₃) ₂ | neat | 1 g |
| DLM-664-1 | 2-Nitroaniline (ring-D ₄ , 98%) | O ₂ NC ₆ D ₄ NH ₂ | neat | 1 g |
| NEW DLM-294-5 | Nitrobenzene (D ₅ , 99%) | C ₆ D ₅ NO ₂ | neat | 5 g |
| NEW DLM-294-10 | | | neat | 10 g |
| ULM-3892-1.2 | Nitrobenzene (unlabeled) | C ₆ H ₅ NO ₂ | 1 mg/mL in acetonitrile | 1.2 mL |
| DLM-295-0.1 | 2-Nitrophenol (ring-D ₄ , 98%) | O ₂ NC ₆ D ₄ OH | neat | 0.1 g |
| DLM-295-0.25 | | | neat | 0.25 g |
| DLM-296-0.1 | 4-Nitrophenol (ring-D ₄ , 98%) | O ₂ NC ₆ D ₄ OH | neat | 0.1 g |
| DLM-296-0.25 | | | neat | 0.25 g |
| ULM-2323-4X25 | n-Nonane (unlabeled) | CH ₃ (CH ₂) ₇ CH ₃ | neat | 4 × 25 mL |
| CLM-6680-1.2 | Octachlorostyrene (¹³ C ₈ , 99%) | *C ₆ Cl ₅ *CCl=*=CCl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| ULM-1709-1.2 | Octachlorostyrene (unlabeled) | C ₆ Cl ₅ CCl=CCl ₂ | 100 µg/mL in isoctane | 1.2 mL |
| OLM-7310-1.2 | Perchloric acid, sodium salt (¹⁸ O ₄ , 90%+) | Cl* ¹⁸ O ₄ ·Na | 100 µg/mL in water | 1.2 mL |
| ULM-7312-1.2 | Perchloric acid, sodium salt (unlabeled) | ClO ₄ ·Na | 100 µg/mL in water | 1.2 mL |
| CLM-216-0.1 | Phenol (¹³ C ₆ , 99%) | *C ₆ H ₅ OH | neat | 0.1 g |
| DLM-695-1 | Phenol (ring-D ₅ , 98%) | C ₆ D ₅ OH | neat | 1 g |
| DLM-695-5 | | | neat | 5g |
| DLM-370-5 | Phenol (D ₆ , 98%) | C ₆ D ₅ OD | neat | 5 g |
| DLM-3039-1MG | Phenylbutazone (diphenyl-D ₁₀ , 98%) | C ₁₉ D ₁₀ H ₁₀ N ₂ O ₂ | neat | 1 mg |
| DLM-3039-0.05 | | | | 0.05 g |
| DLM-3039-0.1 | | | | 0.1 g |
| NEW ULM-7378-1MG | Phenylbutazone (unlabeled) | C ₁₉ H ₂₀ N ₂ O ₂ | neat | 1 mg |
| CLM-3733-1.2 | o-Phenylphenol (ring- ¹³ C ₆ , 99%) | *C ₆ H ₅ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| ULM-7396-1.2 | o-Phenylphenol (unlabeled) | C ₁₂ H ₉ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-3748-1.2 | p-Phenylphenol (ring- ¹³ C ₆ , 99%) CP 96% | *C ₆ H ₅ C ₆ H ₄ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-3040-0.5 | Phthalic acid (carboxyl- ¹³ C, 99%) | C ₆ H ₄ (*CO ₂ H)CO ₂ H | neat | 0.5 g |
| DLM-787-5 | Phthalic acid (ring-D ₄ , 98%) | C ₆ D ₄ (CO ₂ H) ₂ | neat | 5 g |
| DLM-1293-0.1 | 2-Picoline (2-methylpyridine) (D ₇ , 98%) | C ₅ D ₄ NCD ₃ | neat | 0.1 g |
| DLM-1293-1 | | | neat | 1 g |
| DLM-1541-1 | 3-Picoline (3-methylpyridine) (D ₇ , 98%) | C ₅ D ₄ NCD ₃ | neat | 1 g |
| DLM-1294-1 | 4-Picoline (4-methylpyridine) (D ₇ , 98%) | C ₅ D ₄ NCD ₃ | neat | 1 g |
| DLM-1067-5 | 1,2-Propylene oxide (D ₆ , 98%) * | CD ₃ CD ₂ O | neat | 5 g |
| DLM-1158-0.1 | Quinoline (D ₇ , 98%) | C ₉ D ₇ N | neat | 0.1 g |
| DLM-1158-1 | | | neat | 1 g |
| DLM-3322-0.5 | trans-Stilbene (D ₁₂ , 98%) | C ₆ D ₅ CD=CDC ₆ D ₅ | neat | 0.5 g |
| DLM-1083-5 | Styrene (vinyl-D ₃ , 98%) (stabilized with BHT) | C ₆ H ₅ CD=CD ₂ | neat | 5 g |
| DLM-809-5 | Styrene (ring-D ₅ , 98%) (stabilized with BHT) | C ₆ D ₅ CH=CH ₂ | neat | 5 g |
| DLM-380-1.2 | Styrene (D ₈ , 98%) (stabilized with BHT) | C ₆ D ₅ CD=CD ₂ | 100 µg/mL in nonane | 1.2 mL |
| DLM-380-1 | | | neat | 1 g |
| DLM-380-5 | | | neat | 5 g |
| DLM-1088-1 | Terephthalic acid (ring-D ₄ , 98%) | C ₆ D ₄ (CO ₂ H) ₂ | neat | 1 g |
| DLM-1088-5 | | | neat | 5 g |

*Gases require a breakseal flask or cylinder and valve at an additional charge. Breakseal flasks are only available for certain gases at atmospheric pressure.

Priority Pollutant Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------|--|---|--|-----------------------|
| DLM-450-1 | <i>o</i> -Terphenyl (D ₁₄ , 98%) | C ₁₈ D ₁₄ | neat | 1 g |
| DLM-450-5 | | | neat | 5 g |
| DLM-382-1.2 | <i>p</i> -Terphenyl (D ₁₄ , 98%) | C ₁₈ D ₁₄ | 200 µg/mL in isoctane | 1.2 mL |
| DLM-382-1 | | | neat | 1 g |
| DLM-382-5 | | | neat | 5 g |
| ULM-7428-1.2 | <i>p</i> -Terphenyl (unlabeled) | C ₁₈ H ₁₄ | 200 µg/mL in isoctane | 1.2 mL |
| DLM-2279-0.1 | α-Terpineol (propyl methyl-D ₃ , 98%) | CD ₃ C ₆ H ₇ C ₃ H ₇ OH | neat | 0.1 g |
| DLM-2279-0.5 | | | neat | 0.5 g |
| ULM-1704-0.1 | 3,4,5,6-Tetrachlorocatechol (unlabeled) | Cl ₄ C ₆ (OH) ₂ | neat | 0.1 g |
| DLM-35-5 | 1,1,2,2-Tetrachloroethane (D ₂ , 99.6%) | Cl ₂ CDCCl ₂ | neat | 5 g |
| CLM-1965-0.1 | Tetrachloroethylene (¹³ C ₂ , 99%) | Cl ₂ *C=*=CCl ₂ | neat | 0.1 g |
| ULM-1708-0.1 | 3,4,5,6-Tetrachloroguaiacon (unlabeled) | Cl ₄ C ₆ (OH)(OCH ₃) | neat | 0.1 g |
| NEW | ULM-8984-1.2 | Tetrachloro- <i>m</i> -xylene (unlabeled) | C ₈ H ₆ Cl ₄ | 100 µg/mL in isoctane |
| | DLM-2053-0.1 | <i>cis</i> -1,2,3,6-Tetrahydrophthalic anhydride (3,3,4,5,6,6-D ₆ , 98%) | C ₈ D ₆ H ₂ O ₃ | neat |
| DLM-2054-0.1 | <i>cis</i> -1,2,3,6-Tetrahydrophthalimide (3,3,4,5,6,6-D ₆ , 98%) | C ₈ D ₆ H ₃ NO ₂ | neat | 0.1 g |
| CLM-6069-0.1 | Toluene (ring- ¹³ C ₆ , 99%) | *C ₆ H ₅ CH ₃ | neat | 0.1 g |
| CLM-309-0.5 | Toluene (methyl- ¹³ C, 99%) | C ₆ H ₅ *CH ₃ | neat | 0.5 g |
| CLM-309-1 | | | neat | 1 g |
| DLM-1175-1 | Toluene (methyl-D ₃ , 98%) | C ₆ H ₅ CD ₃ | neat | 1 g |
| DLM-1175-5 | | | neat | 5 g |
| DLM-1176-1 | Toluene (ring-D ₅ , 98%) | C ₆ D ₅ CH ₃ | neat | 1 g |
| DLM-1176-5 | | | neat | 5 g |
| DLM-5-5 | Toluene (D ₈ , 99.5%) | C ₆ D ₅ CD ₃ | neat | 5 g |
| DLM-7136-1.2 | Tributyltin chloride (D ₂₇ , 98%) | C ₁₂ D ₂₇ Cl ₃ Sn | 100 µg/mL in MeCl-D ₂ | 1.2 mL |
| ULM-8061-1.2 | Tributyltin chloride (unlabeled) | C ₁₂ H ₂₇ Cl ₃ Sn | 100 µg/mL in MeCl | 1.2 mL |
| ULM-1703-0.1 | 3,4,5-Trichlorocatechol (unlabeled) | Cl ₃ C ₆ H(OH) ₂ | neat | 0.1 g |
| NEW | ULM-9279 | 3,4,6-Trichlorocatechol (unlabeled) | C ₆ H ₃ Cl ₃ O ₂ | Inquire |
| | DLM-1974-0.1 | 1,1,1-Trichloroethane (D ₃ , 98%) | CD ₃ CCl ₃ | neat |
| DLM-1974-1 | | | neat | 0.1 g |
| | | | | 1 g |
| CLM-2075-0.1 | 1,1,2-Trichloroethane (¹³ C ₂ , 99%) | Cl ₂ *CH*CH ₂ Cl | neat | 0.1 g |
| DLM-1975-0.1 | 1,1,2-Trichloroethane (D ₃ , 98%) | Cl ₂ CDCD ₂ Cl | neat | 0.1 g |
| DLM-1975-0.5 | | | neat | 0.5 g |
| CLM-129-0.1 | Trichloroethylene (¹³ C ₂ , 99%) (stabilized with diisopropylamine) | Cl ₂ *C=*=CHCl | neat | 0.1 g |
| DLM-3049-1 | Trichloroethylene (D, 98%) | Cl ₂ C=CDCl | neat | 1 g |
| DLM-2080-0.1 | 1,2,3-Trichloropropane (D ₅ , 98%) CP 95% | CD ₂ ClCDClCD ₂ Cl | neat | 0.1 g |
| DLM-7663 | Triethanolamine (D ₁₅ , 98%) (contains 2-amino-1-propanol) CP 97% | (DOCD ₂ CD ₂) ₃ N | | Inquire |
| DLM-3344-5 | Vinyl bromide (D ₃ , 98%) * (inhibited with hydroquinone) | CD ₂ =CDBr | neat | 5 g |
| DLM-167-1.2 | Vinyl chloride (D ₃ , 98%) | CD ₂ =CDCl | 50 µg/mL in methanol-OD | 1.2 mL |
| DLM-167-5 | Vinyl chloride (D ₃ , 98%) * (inhibited with hydroquinone) | | neat | 5 g |
| DLM-2398-5 | <i>m</i> -Xylene (D ₁₀ , 98%) | C ₆ D ₄ (CD ₃) ₂ | neat | 5 g |
| DLM-808-5 | <i>o</i> -Xylene (D ₁₀ , 98%) | C ₆ D ₄ (CD ₃) ₂ | neat | 5 g |
| DLM-313-5 | <i>p</i> -Xylene (D ₁₀ , 98%) | C ₆ D ₄ (CD ₃) ₂ | neat | 5 g |

Notes

A photograph showing two young boys from behind, standing in a body of water heavily polluted with trash. The water is dark and filled with plastic bottles, cans, and other debris. In the background, there's a large industrial building with rusted pipes and a fence. The boy on the left wears a white t-shirt and light blue shorts, while the boy on the right wears a dark blue t-shirt and grey cargo shorts.

Chlorinated benzenes and phenols are common chemical contaminants produced in vast quantities and used in many industrial processes, bringing them in contact with the environment in a variety of locations. Owing to the sheer volume of use in commercial products, halogenated benzenes and phenols remain among the largest contributors to environmental contamination.

Chlorobenzene and Chlorophenol Standard Mixtures

Owing to the sheer volume of use in commercial products, halogenated benzenes and phenols remain among the largest contributors to environmental contamination. The standards listed here are in routine use in many laboratories around the world.

US EPA Method 1653A

US EPA Method 1653A is used for the determination of pollutants in pulp and paper industry wastewater. This revision was promulgated in 1997, superseding the earlier method 1653. While still used primarily for the determination of chlorophenolic compounds, Revision A incorporates several changes to the analytical procedure, including the use of specially formulated standard mixtures applicable to this revision.

US EPA CLP DMC Standard Mixtures

EPA's Contract Laboratory Program (CLP) has developed methods for the analysis of volatile and semi-volatile compounds which utilize isotopically labeled internal standards. These deuterated monitoring compounds (DMCs) have been added to strengthen the analysis by providing sample-by-sample internal standard addition. CIL's CLP DMC standard mixtures are designed to match requirements of the OLC and SOM test methods.

US EPA Method 1624/1625

CIL maintains a full suite of standards used for the analysis of volatile and semi-volatile organic compounds by US EPA Method 1624/1625.

Human Exposure Analysis

Exposure analysis studies the ultimate effects environmental or food pollutants have on a biological system. Increasingly, public health studies are looking at the relationship between adverse health issues and exposure to environmental and food pollutants, as well as consumption pollutants such as tobacco compounds, bisphenols, etc. CIL has worked for years to develop analytical standards to assist exposure research and now offers some targeted mixtures as well. Our "phenolics" mixes include several ubiquitous exposure biomarkers, including PCB metabolites, BDE metabolites, tetrabromobisphenol A, triclosan, chlorophenols, and bromophenols. The analytes in the calibration solutions are the derivatized analogues of the target compounds, and the unusual concentrations are a result of corrections for molecular weights of the derivatized products. The spiking solution contains underderivatized compounds to allow for derivatization during the isotope dilution method.

Chlorobenzene and Chlorophenol Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|--------------------|
| EM-1724-A | ¹³ C-Labeled Chlorobenzene Cocktail Solution – Mono, Di, Tri Isomers | 1.2 mL in isoctane |
| EM-1724-B | ¹³ C-Labeled Chlorobenzene Cocktail Solution – Mono, Di, Tri Isomers | 1.2 mL in methanol |
| | Labeled | (μ g/mL) |
| | Chlorobenzene (¹³ C ₆ , 99%) | 100 |
| | 1,4-Dichlorobenzene (¹³ C ₆ , 99%) | 100 |
| | 1,2,4-Trichlorobenzene (¹³ C ₆ , 99%) | 100 |
| EM-1725-A | ¹³ C-Labeled Chlorobenzene Cocktail Solution – Tetra, Penta, Hexa Isomers | 1.2 mL in isoctane |
| EM-1725-B | ¹³ C-Labeled Chlorobenzene Cocktail Solution – Tetra, Penta, Hexa Isomers | 1.2 mL in methanol |
| | Labeled | (μ g/mL) |
| | 1,2,4,5-Tetrachlorobenzene (¹³ C ₆ , 99%) | 100 |
| | Pentachlorobenzene (¹³ C ₆ , 99%) | 100 |
| | Hexachlorobenzene (¹³ C ₆ , 99%) | 100 |
| EM-1726-A | ¹³ C-Labeled Chlorophenol Cocktail Solution – Mono, Di, Tri Isomers | 1.2 mL in isoctane |
| EM-1726-B | ¹³ C-Labeled Chlorophenol Cocktail Solution – Mono, Di, Tri Isomers | 1.2 mL in methanol |
| | Labeled | (μ g/mL) |
| | 4-Chlorophenol (¹³ C ₆ , 99%) | 100 |
| | 2,4-Dichlorophenol (¹³ C ₆ , 99%) | 100 |
| | 2,4,6-Trichlorophenol (¹³ C ₆ , 99%) | 100 |
| EM-1727-A | ¹³ C-Labeled Chlorophenol Cocktail Solution – Tri, Tetra, Penta Isomers | 1.2 mL in isoctane |
| EM-1727-B | ¹³ C-Labeled Chlorophenol Cocktail Solution – Tri, Tetra, Penta Isomers | 1.2 mL in methanol |
| | Labeled | (μ g/mL) |
| | 2,4,5-Trichlorophenol (¹³ C ₆ , 99%) | 100 |
| | 2,3,4,5-Tetrachlorophenol (¹³ C ₆ , 99%) | 100 |
| | Pentachlorophenol (¹³ C ₆ , 99%) | 100 |
| ES-5401 | ¹³ C-Labeled Mono-Hexa Chlorobenzene Solution | 1.2 mL in toluene |
| | Labeled | (μ g/mL) |
| | Chlorobenzene (¹³ C ₆ , 99%) | 500 |
| | 1,4-Dichlorobenzene (¹³ C ₆ , 99%) | 500 |
| | 1,2,3-Trichlorobenzene (¹³ C ₆ , 99%) | 500 |
| | 1,2,3,4-Tetrachlorobenzene (¹³ C ₆ , 99%) | 500 |
| | Pentachlorobenzene (¹³ C ₆ , 99%) | 500 |
| | Hexachlorobenzene (¹³ C ₆ , 99%) | 500 |
| ES-5406 | Native Mono-Hexa Chlorobenzene Solution | 1.2 mL in toluene |
| | Unlabeled | (μ g/mL) |
| | Chlorobenzene | 500 |
| | 1,4-Dichlorobenzene | 500 |
| | 1,2,3-Trichlorobenzene | 500 |
| | 1,2,3,4-Tetrachlorobenzene | 500 |
| | Pentachlorobenzene | 500 |
| | Hexachlorobenzene | 500 |

US EPA Method 1653A Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|------------------|
| EM-4173 | Method 1653A Labeled Chlorophenolic Derivatives Mixture – 1 Each: EM-4173-1 and EM-4173-2 (Note: unlabeled internal standard, 3,4,5-trichlorophenol already formulated into the standard) | 2 x 1 mL |
| EM-4173-1 | Method 1653A Labeled Chlorophenolic Derivatives Mixture | 1 mL in methanol |
| EM-4173-2 | Method 1653A Labeled Chlorophenolic Derivatives Mixture | 1 mL in acetone |

| Labeled | EM-4173-1 (μ g/mL) | EM-4173-2 (μ g/mL) |
|---|----------------------------|----------------------------|
| 2,4-Dichlorophenol ($^{13}\text{C}_6$, 99%) | 250 | – |
| 4-Chloroguaiacol ($^{13}\text{C}_6$, 99%) | 250 | – |
| 4,5-Dichlorocatechol (ring- $^{13}\text{C}_6$, 99%) | 250 | – |
| 4,5,6-Trichloroguaiacol ($^{13}\text{C}_6$, 99%) | 250 | – |
| Pentachlorophenol ($^{13}\text{C}_6$, 99%) | 250 | – |
| 3,4,5,6-Tetrachloroguaiacol ($^{13}\text{C}_6$, 99%) | 250 | – |
| 3,4,5,6-Tetrachlorocatechol (ring- $^{13}\text{C}_6$, 99%) | 250 | – |
| 3,4,5-Trichlorophenol (unlabeled) (IS) | 250 | – |
| 5-Chlorovanillin (ring- $^{13}\text{C}_6$, 99%) | – | 250 |

| | | |
|---------|--|------------------|
| EM-4181 | Method 1653A Regulated Chlorophenolics Mixture-1 | 1 mL in methanol |
|---------|--|------------------|

| Unlabeled | (μ g/mL) |
|-----------------------------|---------------|
| Pentachlorophenol | 1000 |
| 3,4,5,6-Tetrachloroguaiacol | 1000 |
| 3,4,5-Trichlorocatechol | 1000 |
| 3,4,5-Trichloroguaiacol | 500 |
| 4,5,6-Trichloroguaiacol | 500 |
| 2,4,6-Trichlorophenol | 500 |
| 3,4,5,6-Tetrachlorocatechol | 1000 |
| 2,3,4,6-Tetrachlorophenol | 500 |
| 3,4,6-Trichlorocatechol | 1000 |
| 3,4,6-Trichloroguaiacol | 500 |
| 2,4,5-Trichlorophenol | 500 |

| | | |
|---------|--|-----------------|
| EM-4182 | Method 1653A Regulated Chlorophenolics Mixture-2 | 1 mL in acetone |
|---------|--|-----------------|

| Unlabeled | (μ g/mL) |
|-------------------|---------------|
| Trichlorosyringol | 500 |

| | | |
|---------|--|------------------|
| EM-4183 | Method 1653A Other Chlorophenolics Mixture-1 | 1 mL in methanol |
|---------|--|------------------|

| Unlabeled | (μ g/mL) |
|----------------------|---------------|
| 4-Chlorocatechol | 250 |
| 4-Chlorophenol | 250 |
| 3,6-Dichlorocatechol | 500 |
| 3,4-Dichloroguaiacol | 500 |
| 4,6-Dichloroguaiacol | 500 |
| 2,6-Dichlorophenol | 500 |
| 4-Chloroguaiacol | 250 |
| 3,4-Dichlorocatechol | 500 |
| 4,5-Dichlorocatechol | 500 |
| 4,5-Dichloroguaiacol | 500 |
| 2,4-Dichlorophenol | 500 |

US EPA Method 1653A Standard Mixtures

| Catalog No. | Compound | Amount |
|---------------------------------------|--|-----------------|
| EM-4184 | Method 1653A Other Chlorophenolics Mixture-2 | 1 mL in acetone |
| Unlabeled ($\mu\text{g/mL}$) | | |
| | 2-Chlorosyringealdehyde | 500 |
| | 5-Chlorovanillin | 500 |
| | 6-Chlorovanillin | 500 |
| | 2,6-Dichlorosyringealdehyde | 1000 |
| | 5,6-Dichlorovanillin | 1000 |
| EM-4185 | Set of Regulated Chlorophenolics Mixtures 1 Each: EM-4181 and EM-4182 | 2 \times 1 mL |
| EM-4186 | Set of Other Chlorophenolics Mixtures 1 Each: EM-4183 and EM-4184 | 2 \times 1 mL |
| EM-4180 | Set of Chlorophenolics Mixtures 1 Each: EM-4181, EM-4182, EM-4183 and EM-4184 | 4 \times 1 mL |

US EPA Method 1653 Standard Mixtures

| | | |
|---------------------------------------|---|-----------------|
| EM-4018 | Method 1653 Unlabeled Chloroguaiacol Cocktail | 1 mL in acetone |
| Unlabeled ($\mu\text{g/mL}$) | | |
| | 4-Chloroguaiacol | 250 |
| | 3,4-Dichloroguaiacol | 500 |
| | 4,5-Dichloroguaiacol | 500 |
| | 4,6-Dichloroguaiacol | 500 |
| | 3,4,5-Trichloroguaiacol | 500 |
| | 3,4,6-Trichloroguaiacol | 500 |
| | 4,5,6-Trichloroguaiacol | 500 |
| | 3,4,5,6-Tetrachloroguaiacol | 1000 |
| EM-4028 | Instrument Performance Standard | 1 mL in acetone |
| Unlabeled ($\mu\text{g/mL}$) | | |
| | 2,2'-Difluorobiphenyl | 5000 |

US EPA CLP DMC Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---------------------------------------|-------------------------------|
| ES-5037 | CLP Semi-Volatiles DMC Stock Solution | 1.2 mL in MeCl-D ₂ |

| Labeled | (μ g/mL) |
|---|---------------|
| Phenol (ring-D ₅ , 98%) | 2000 |
| Bis(2-chloroethyl) ether (D ₈ , 98%) | 2000 |
| 2-Chlorophenol (ring-D ₄ , 98%) | 2000 |
| 4-Methylphenol (<i>p</i> -Cresol) (D ₈ , 98%) | 2000 |
| Nitrobenzene (D ₅ , 99%) | 2000 |
| 2-Nitrophenol (ring-D ₄ , 98%) | 2000 |
| 2,4-Dichlorophenol (ring-D ₃ , 98%) | 2000 |
| 4-Chloroaniline (D ₄ , 98%) | 2000 |
| Dimethyl phthalate (dimethyl-D ₆ , 98%) | 2000 |
| Acenaphthylene (D ₈ , 98%) | 2000 |
| 4-Nitrophenol (ring-D ₄ , 98%) | 2000 |
| Fluorene (D ₁₀ , 98%) | 2000 |
| 4,6-Dinitro-2-methylphenol (ring-D ₂ , 98%) | 2000 |
| Anthracene (D ₁₀ , 98%) | 2000 |
| Pyrene (D ₁₀ , 98%) | 2000 |
| Benzo[a]pyrene (D ₁₂ , 98%) | 2000 |

| | | |
|-----------|--|-----------------------|
| ES-5038 | CLP OLC Volatiles DMC Stock Solutions 1 Each: ES-5038-1 and ES-5038-2 | 1 set |
| ES-5038-1 | CLP OLC Volatiles Non-Ketone DMC Stock Solution | 1 mL in methanol-OD |
| ES-5038-2 | CLP OLC Volatiles Ketone DMC Stock Solution | 0.5 mL in methanol-OD |

| Labeled | ES-5038-1 (μ g/mL) | ES-5038-2 (μ g/mL) |
|--|----------------------------|----------------------------|
| Vinyl chloride (D ₃ , 98%) | 100 | — |
| Chloroethane (D ₅ , 98%) | 100 | — |
| 1,1-Dichloroethene (D ₂ , 98%) | 100 | — |
| 2-Butanone (1,1,1,3-D ₅ , 98%) | — | 200 |
| Chloroform (D, 98%) | 100 | — |
| 1,2-Dichloroethane (D ₄ , 99%) | 100 | — |
| Benzene (D ₆ , 99.5%) | 100 | — |
| 1,2-Dichloropropane (D ₆ , 98%) | 100 | — |
| Toluene (D ₈ , 99.5%) | 100 | — |
| 1,3-Dichloropropene (D ₄ , 98%) (<i>cis/trans</i> mixture) | 100 | — |
| 2-Hexanone (1,1,1,3-D ₅ , 98%) | — | 200 |
| Bromoform (D, 99.5%) | 100 | — |
| 1,1,2,2-Tetrachloroethane (D ₂ , 99.6%) | 100 | — |
| 1,2-Dichlorobenzene (D ₄ , 99%) | 100 | — |

US EPA CLP DMC Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|---------------------|
| ES-5286 | CLP SOM Volatiles Non-Ketone DMC Stock Solution | 1 mL in methanol-OD |
| ES-5286-10X | CLP SOM Volatiles Non-Ketone DMC 10X Stock Solution | 1 mL in methanol-OD |

| Labeled | ES-5286 (μ g/mL) | ES-5286-10X (μ g/mL) |
|--|--------------------------|------------------------------|
| Vinyl chloride (D_3 , 98%) | 100 | 1000 |
| Chloroethane (D_5 , 98%) | 100 | 1000 |
| 1,1-Dichloroethylene (2,2- D_2 , 98%) | 100 | 1000 |
| Chloroform (D, 99.8%) | 100 | 1000 |
| 1,2-Dichloroethane (D_4 , 99%) | 100 | 1000 |
| Benzene (D_6 , 99.5%) | 100 | 1000 |
| 1,2-Dichloropropane (D_6 , 98%) | 100 | 1000 |
| Toluene (D_8 , 99.5%) | 100 | 1000 |
| 1,3-Dichloropropene (D_4 , 98%) (<i>cis/trans</i> mixture) | 100 | 1000 |
| 1,1,2,2-Tetrachloroethane (D_2 , 99.6%) | 100 | 1000 |
| 1,2-Dichlorobenzene (D_4 , 99%) | 100 | 1000 |

| | | |
|-------------|---|-----------------------|
| ES-5287 | CLP SOM Volatiles Ketone DMC Stock Solution | 0.5 mL in methanol-OD |
| ES-5287-10X | CLP SOM Volatiles Ketone DMC 10X Stock Solution | 0.5 mL in methanol-OD |

| Labeled | ES-5287 (μ g/mL) | ES-5287-10X (μ g/mL) |
|-------------------------------------|--------------------------|------------------------------|
| 2-Butanone (1,1,1,3,3- D_5 , 98%) | 500 | 5000 |
| 2-Hexanone (1,1,1,3,3- D_5 , 98%) | 500 | 5000 |

| | | |
|-------------|--|---------------------|
| ES-5288 | CLP SOM Volatiles 1,4-Dioxane DMC Stock Solution | 1 mL in methanol-OD |
| ES-5288-10X | CLP SOM Volatiles 1,4-Dioxane 10X DMC Stock Solution | 1 mL in methanol-OD |

| Labeled | ES-5288 (μ g/mL) | ES-5288-10X (μ g/mL) |
|---|--------------------------|------------------------------|
| 1,4-Dioxane (<i>p</i> -dioxane) (D_8 , 99%) | 1250 | 12,500 |

US EPA Methods 1624/1625 Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|--|---|
| ES-2036 | Acid Extractables Mixture-3 | 1 mL in benzene-D ₆ |
| | Labeled | (μ g/mL) |
| | (EPA 222A) 4-Chloro-3-methylphenol (ring-2,6-D ₂ , 98%) | 5000 |
| | (EPA 224A) 2-Chlorophenol (ring-D ₄ , 98%) | 5000 |
| | (EPA 231A) 2,4-Dichlorophenol (ring-D ₃ , 98%) | 5000 |
| | (EPA 234A) 2,4-Dimethylphenol (ring-D ₃ , 98%) | 5000 |
| | (EPA 260A) 4,6-Dinitro-2-methylphenol (ring-D ₂ , 98%) | 5000 |
| | (EPA 259A) 2,4-Dinitrophenol (ring-D ₃ , 98%) | 5000 |
| | (EPA 257A) 2-Nitrophenol (ring-D ₄ , 98%) | 5000 |
| | (EPA 258A) 4-Nitrophenol (ring-D ₄ , 98%) | 5000 |
| | (EPA 264A) Pentachlorophenol (¹³ C ₆ , 99%) | 5000 |
| | (EPA 265A) Phenol (ring-D ₅ , 98%) | 5000 |
| | (EPA 631A) 2,4,5-Trichlorophenol (ring-D ₂ , 98%) | 5000 |
| | (EPA 221A) 2,4,6-Trichlorophenol (ring-D ₂ , 98%) | 5000 |
| ES-2002 | Base Neutrals Mixture-4.3 | 1 mL in benzene-D ₆ |
| | Labeled | (μ g/mL) |
| | (EPA 241B) 4-Bromophenyl phenyl ether (phenyl-D ₅ , 98%) | 5000 |
| | (EPA 220B) 2-Chloronaphthalene (D ₇ , 98%) | 5000 |
| | (EPA 240B) 4-Chlorophenyl phenyl ether (phenyl-D ₅ , 98%) | 5000 |
| | (EPA 268B) Di-n-butyl phthalate (ring-D ₄ , 98%) | 5000 |
| | (EPA 270B) Diethyl phthalate (ring-D ₄ , 98%) | 5000 |
| | (EPA 269B) Di-n-octyl phthalate (ring-D ₄ , 98%) | 5000 |
| | (EPA 209B) Hexachlorobenzene (¹³ C ₆ , 99%) | 5000 |
| | (EPA 212B) Hexachloroethane (1- ¹³ C, 99%) | 5000 |
| | (EPA 254B) Isophorone (3-methyl-D ₃ ; 2,4,4,6,6-D ₅ , 98%) | 5000 |
| | (EPA 208B) 1,2,4-Trichlorobenzene (D ₃ , 98%) | 5000 |
| ES-2026 | Base Neutrals Dilution Mixture-5.2 | 1 mL in benzene-D ₆ |
| | Labeled | (μ g/mL) |
| | (EPA 628B) Carbazole (ring-D ₈ , 98%) | 5000 |
| | (EPA 261B) N-Nitrosodimethylamine (D ₆ , 98%) | 5000 |
| | (EPA 263B) N-Nitrosodi-n-propylamine (D ₁₄ , 98%) | 5000 |
| ES-2003 | Base Neutrals Mixture-6.2 | 2 × 1 mL in 50% benzene-D ₆ and 50% MeCl-D ₂ |
| | Labeled | (μ g/mL) |
| | (EPA 201B) Acenaphthene (D ₁₀ , 99%) | 2500 |
| | (EPA 278B) Anthracene (D ₁₀ , 98%) | 2500 |
| | (EPA 275B) Benzo[k]fluoranthene (D ₁₂ , 98%) | 2500 |
| | (EPA 218B) Bis(2-chloroethyl) ether (D ₈ , 98%) | 2500 |
| | (EPA 276B) Chrysene (D ₁₂ , 98%) | 2500 |
| | (EPA 280B) Fluorene (D ₁₀ , 98%) | 2500 |
| | (EPA 255B) Naphthalene (D ₈ , 99%) | 2500 |
| | (EPA 284B) Pyrene (D ₁₀ , 98%) | 2500 |

US EPA Methods 1624/1625 Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------|--|---------------------------------|
| ES-2004 | Base Neutrals Mixture-6.3 | 1 mL in benzene-D ₆ |
| Labeled | | |
| (EPA 272B) | Benz[a]anthracene (D ₁₂ , 98%) | 5000 |
| (EPA 267B) | Butyl benzyl phthalate (ring-D ₄ , 99%) | 5000 |
| (EPA 243B) | Bis(2-chloroethoxy) methane (chloroethoxy-D ₈ , 98%) | 5000 |
| (EPA 266B) | Bis(2-ethylhexyl) phthalate (ring-D ₄ , 99%) | 5000 |
| (EPA 282B) | Dibenz[a,h]anthracene (D ₁₄ , 98%) | 5000 |
| (EPA 225B) | 1,2-Dichlorobenzene (D ₄ , 99%) | 5000 |
| (EPA 227B) | 1,4-Dichlorobenzene (D ₄ , 98%) | 5000 |
| (EPA 271B) | Dimethyl phthalate (ring-D ₄ , 99%) | 5000 |
| (EPA 236B) | 2,6-Dinitrotoluene (methyl-D ₃ , 98%) | 5000 |
| (EPA 256B) | Nitrobenzene (D ₅ , 99%) | 5000 |
| (EPA 629B) | 1,2,3-Trichlorobenzene (D ₃ , 98%) | 5000 |
| <hr/> | | |
| ES-2032 | Purgeables/Volatiles Mixture-E.1 | 1 mL in methanol-D ₄ |
| Labeled | | |
| (EPA 206V) | Carbon tetrachloride (¹³ C, 99%) | 50 |
| (EPA 207V) | Chlorobenzene (D ₅ , 99%) | 50 |
| (EPA 223V) | Chloroform (¹³ C, 99%) | 50 |
| (EPA 213V) | 1,1-Dichloroethane (2,2,2-D ₃ , 98%) | 50 |
| (EPA 229V) | 1,1-Dichloroethylene (2,2-D ₂ , 98%) | 50 |
| (EPA 244V) | Methylene chloride (D ₂ , 99.9%) | 50 |
| (EPA 232V) | 1,2-Dichloropropane (D ₆ , 98%) | 50 |
| (EPA 214V) | 1,1,2-Trichloroethane (¹³ C ₂ , 99%) | 50 |
| <hr/> | | |
| ES-2006 | Purgeables/Volatiles Mixture-F | 1 mL in methanol-D ₄ |
| Labeled | | |
| (EPA 204V) | Benzene (D ₆ , 99.5%) | 50 |
| (EPA 247V) | Bromoform (¹³ C, 99%) | 50 |
| (EPA 210V) | 1,2-Dichloroethane (D ₄ , 99%) | 50 |
| (EPA 238V) | Ethylbenzene (D ₁₀ , 98%) | 50 |
| (EPA 215V) | 1,1,2,2-Tetrachloroethane (D ₂ , 99.6%) | 50 |
| (EPA 286V) | Toluene (D ₈ , 99.5%) | 50 |
| (EPA 211V) | 1,1,1-Trichloroethane (D ₃ , 98%) | 50 |
| <hr/> | | |
| ES-2008 | Purgeables/Volatiles Mixture-H | 1 mL in methanol-D ₄ |
| Labeled | | |
| (EPA 248V) | Bromodichloromethane (¹³ C, 99%) | 50 |
| (EPA 251V) | Chlorodibromomethane (¹³ C, 99%) | 50 |
| (EPA 30V) | 1,2-Dichloroethylene (1,2-D ₂ , 98%) | 50 |
| (EPA 33V) | 1,3-Dichloropropene (D ₄ , 98%) (<i>cis/trans</i> mixture) | 50 |
| (EPA 627V) | 1,4-Dioxane (D ₈ , 99%) | 50 |
| (EPA 285V) | Tetrachloroethylene (¹³ C ₂ , 99%) | 50 |
| (EPA 287V) | Trichloroethylene (¹³ C ₂ , 99%) | 50 |

Human Exposure Analysis Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|---|---|
| NEW ES-5510 | Extended Phenolic Calibration Solutions [CS1-CS6] | 6 x 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS1 | Extended Phenolic Calibration Solution [CS1] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS2 | Extended Phenolic Calibration Solution [CS2] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS3 | Extended Phenolic Calibration Solution [CS3] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS4 | Extended Phenolic Calibration Solution [CS4] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS5 | Extended Phenolic Calibration Solution [CS5] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |
| NEW ES-5510-CS6 | Extended Phenolic Calibration Solution [CS6] | 0.5 mL in 80% nonane/20% dodecane, hexane and toluene |

All concentrations are in ng/mL (ppb)*

| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
|--|-------|------|------|------|------|------|------|
| 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl | | 0.21 | 0.52 | 2.60 | 21 | 104 | 520 |
| 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl | | 0.21 | 0.52 | 2.59 | 21 | 104 | 519 |
| 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl | | 0.21 | 0.52 | 2.59 | 26.2 | 103 | 517 |
| 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) | | 2.10 | 5.24 | 26.2 | 210 | 1048 | 5242 |
| 4-Methoxy-2,2',3,4'-tetrabromodiphenyl ether | | 0.21 | 0.51 | 2.57 | 21 | 103 | 514 |
| 3-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | | 0.21 | 0.51 | 2.57 | 21 | 103 | 514 |
| 5-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | | 0.21 | 0.51 | 2.57 | 21 | 103 | 514 |
| 6-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | | 0.21 | 0.51 | 2.57 | 21 | 103 | 514 |
| 4'-Methoxy-2,2',4,5'-tetrabromodiphenyl ether | | 0.21 | 0.51 | 2.57 | 21 | 102 | 514 |
| 5'-Methoxy-2,2',4,4',5-pentabromodiphenyl ether | | 0.20 | 0.51 | 2.56 | 21 | 102 | 512 |
| 6-Methoxy-2,2',4,4',5-pentabromodiphenyl ether | | 0.20 | 0.51 | 2.56 | 21 | 102 | 512 |
| 2,4-Dibromoanisole | | 0.21 | 0.53 | 2.64 | 21 | 106 | 528 |
| 2,4,5-Tribromoanisole | | 0.21 | 0.52 | 2.61 | 21 | 104 | 521 |
| 2,4,6-Tribromoanisole | | 0.21 | 0.52 | 2.61 | 21 | 104 | 521 |
| Pentabromoanisole | | 0.21 | 0.51 | 2.57 | 21 | 103 | 514 |
| Pentachloroanisole | | 2.11 | 5.26 | 26.3 | 210 | 1053 | 5263 |
| Dimethyl tetrabromobisphenol A | | 0.11 | 0.26 | 1.31 | 11 | 53 | 263 |

Labeled

| | | | | | | | |
|--|--|------|------|------|------|------|------|
| 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl (¹³ C ₁₂ , 99%) | | 104 | 104 | 104 | 104 | 104 | 104 |
| 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl (¹³ C ₁₂ , 99%) | | 104 | 104 | 104 | 104 | 104 | 104 |
| 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl (¹³ C ₁₂ , 99%) | | 103 | 103 | 103 | 103 | 103 | 103 |
| 6-Methoxy-2,2',4,4'-tetrabromodiphenyl ether (ring- ¹³ C ₁₂ , 99%) | | 103 | 103 | 103 | 103 | 103 | 103 |
| 6'-Methoxy-2,2',4,4',6-pentabromodiphenyl ether (¹³ C ₁₂ , 99%) | | 102 | 102 | 102 | 102 | 102 | 102 |
| 2,4-Dibromoanisole (ring- ¹³ C ₆ , 99%) | | 105 | 105 | 105 | 105 | 105 | 105 |
| 2,4,5-Tribromoanisole (ring- ¹³ C ₆ , 99%) | | 104 | 104 | 104 | 104 | 104 | 104 |
| 2,4,6-Tribromoanisole (ring- ¹³ C ₆ , 99%) | | 104 | 104 | 104 | 104 | 104 | 104 |
| Pentabromoanisole (¹³ C ₆ , 99%) | | 103 | 103 | 103 | 103 | 103 | 103 |
| Pentachloroanisole (¹³ C ₆ , 99%) | | 1052 | 1052 | 1052 | 1052 | 1052 | 1052 |
| Dimethyl tetrabromobisphenol A (ring- ¹³ C ₁₂ , 99%) | | 53 | 53 | 53 | 53 | 53 | 53 |
| 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) (¹³ C ₁₂ , 99%) | | 1047 | 1047 | 1047 | 1047 | 1047 | 1047 |
| 1,2,3,4-Tetrachlorodibenzo-p-dioxin (¹³ C ₆ , 99%) | | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl (¹³ C ₁₂ , 99%) | | 208 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-Tetrabromodiphenyl ether (¹³ C ₁₂ , 99%) | | 77 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-Hexabromodiphenyl ether (¹³ C ₁₂ , 99%) | | 139 | 75 | 75 | 75 | 75 | 75 |

*Concentrations are determined for listed derivatized compounds.

Human Exposure Analysis Standard Mixtures

| Catalog No. | Compound | Amount |
|-----------------------------|--|---|
| NEW ES-5511 | Extended Phenolic Spiking Solution | 10 mL in methanol |
| | Labeled | (ng/mL) |
| | 4-Hydroxy-2,3,3',4',5-pentachlorobiphenyl (¹³ C ₁₂ , 99%) | 10 |
| | 4-Hydroxy-2,2',3,4',5,5'-hexachlorobiphenyl (¹³ C ₁₂ , 99%) | 10 |
| | 4-Hydroxy-2,2',3,4',5,5',6-heptachlorobiphenyl (¹³ C ₁₂ , 99%) | 10 |
| | 6-Hydroxy-2,2',4,4'-tetrabromodiphenyl ether (¹³ C ₁₂ , 99%) | 10 |
| | 6'-Hydroxy-2,2',4,4',6-pentabromodiphenyl ether (¹³ C ₁₂ , 99%) | 10 |
| | 2,4-Dibromophenol (¹³ C ₆ , 99%) | 10 |
| | 2,4,5-Tribromophenol (¹³ C ₆ , 99%) | 10 |
| | 2,4,6-Tribromophenol (¹³ C ₆ , 99%) | 10 |
| | Pentabromophenol (¹³ C ₆ , 99%) | 10 |
| | Pentachlorophenol (¹³ C ₆ , 99%) | 100 |
| | Tetrabromobisphenol A (ring- ¹³ C ₁₂ , 99%) | 5 |
| | 5-Chloro-2-(2,4-dichlorophenoxy)-phenol (triclosan) (¹³ C ₁₂ , 99%) | 100 |
| ES-5321 | Multi-Analyte Recovery Spiking Standard | 10 mL in 88% hexane/ 2% dodecane/10% nonane |
| NEW ES-5321-200X-1.2 | Multi-Analyte Recovery Spiking Standard | 1.2 mL in nonane |
| | Labeled | ES-5321 ES-5321-200X-1.2 (ng/mL) (ng/mL) |
| | 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 2.5 500 |
| | 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) | 208 10.0 2000 |
| | 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) | 77 7.5 1500 |
| | 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) | 139 7.5 1500 |
| ES-9444 | Extended Phenolic Native PAR Standard | 1.2 mL in 91% nonane/ 9% toluene |
| | Unlabeled | (ng/mL)* |
| | 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl | 520 |
| | 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl | 519 |
| | 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl | 517 |
| | 4-Methoxy-2,2',3,4'-tetrabromodiphenyl ether | 514 |
| | 3-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | 514 |
| | 5-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | 514 |
| | 6-Methoxy-2,2',4,4'-tetrabromodiphenyl ether | 514 |
| | 4'-Methoxy-2,2',4,5'-tetrabromodiphenyl ether | 514 |
| | 5'-Methoxy-2,2',4,4',5-pentabromodiphenyl ether | 512 |
| | 6-Methoxy-2,2',4,4',5-pentabromodiphenyl ether | 512 |
| | 2,4-Dibromoanisole | 528 |
| | 2,4,5-Tribromoanisole | 521 |
| | 2,4,6-Tribromoanisole | 521 |
| | Dimethyl tetrabromobisphenol A | 263 |
| | Pentabromoanisole | 514 |
| | 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) | 5240 |
| | Pentachloroanisole | 5263 |

*Concentrations are determined for listed derivatized compounds.

Human Exposure Analysis Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|--|----------------------|
| NEW ES-5482 | Phenolic Calibration Standards [CS1-CS6] | 6 x 0.5 mL in nonane |
| NEW ES-5482-CS1 | Phenolic Calibration Standard [CS1] | 0.5 mL in nonane |
| NEW ES-5482-CS2 | Phenolic Calibration Standard [CS2] | 0.5 mL in nonane |
| NEW ES-5482-CS3 | Phenolic Calibration Standard [CS3] | 0.5 mL in nonane |
| NEW ES-5482-CS4 | Phenolic Calibration Standard [CS4] | 0.5 mL in nonane |
| NEW ES-5482-CS5 | Phenolic Calibration Standard [CS5] | 0.5 mL in nonane |
| NEW ES-5482-CS6 | Phenolic Calibration Standard [CS6] | 0.5 mL in nonane |

| <i>All concentrations are in ng/mL (ppb)*</i> | | | | | | |
|--|-------|------|------|------|-----|-----|
| Unlabeled | IUPAC | CS1 | CS2 | CS3 | CS4 | CS5 |
| 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl | | 0.52 | 1.04 | 10.4 | 104 | 520 |
| 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl | | 0.52 | 1.04 | 10.4 | 104 | 519 |
| 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl | | 0.52 | 1.03 | 10.3 | 103 | 517 |
| Pentachloroanisole | | 0.53 | 1.05 | 10.5 | 105 | 526 |
| 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) | | 0.52 | 1.05 | 10.5 | 105 | 524 |
| Pentabromoanisole | | 0.51 | 1.03 | 10.3 | 103 | 514 |
| Labeled | | | | | | |
| 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl (¹³ C ₁₂ , 99%) | | 104 | 104 | 104 | 104 | 104 |
| 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl (¹³ C ₁₂ , 99%) | | 104 | 104 | 104 | 104 | 104 |
| 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl (¹³ C ₁₂ , 99%) | | 103 | 103 | 103 | 103 | 103 |
| Pentachloroanisole (¹³ C ₆ , 99%) | | 263 | 263 | 263 | 263 | 263 |
| 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) (ring- ¹³ C ₁₂ , 99%) | | 262 | 262 | 262 | 262 | 262 |
| Pentabromoanisole (¹³ C ₆ , 99%) | | 103 | 103 | 103 | 103 | 103 |
| 1,2,3,4-Tetrachlorodibenzo-p-dioxin (¹³ C ₆ , 99%) | | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6'-Nonachlorobiphenyl (¹³ C ₁₂ , 99%) | 208 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-Tetrabromodiphenyl ether (¹³ C ₁₂ , 99%) | 77 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-Hexabromodiphenyl ether (¹³ C ₁₂ , 99%) | 139 | 75 | 75 | 75 | 75 | 75 |

*Concentrations are determined for listed derivatized compounds.

Human Exposure Analysis Standard Mixtures

| Catalog No. | Compound | Amount | |
|---|---|-----------------------------|------|
| NEW ES-5483 | Phenolic Spiking Standard | 10 mL in acetonitrile | |
| Labeled | | | |
| | 4-Hydroxy-2,3,3',4',5-pentachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 10 | |
| | 4-Hydroxy-2,2',3,4',5,5'-hexachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 10 | |
| | 4-Hydroxy-2,2',3,4',5,5',6-heptachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 10 | |
| | Pentachlorophenol ($^{13}\text{C}_6$, 99%) | 25 | |
| | 5-Chloro-2-(2,4-dichlorophenoxy)-phenol (triclosan) ($^{13}\text{C}_{12}$, 99%) | 25 | |
| | Pentabromophenol ($^{13}\text{C}_6$, 99%) | 10 | |
| ES-5321 Multi-Analyte Recovery Spiking Standard | | | |
| NEW ES-5321-200X-1.2 Multi-Analyte Recovery Spiking Standard | | | |
| | | | |
| Labeled | | ES-5321 (ng/mL) | |
| | IUPAC | ES-5321-200X-1.2 (ng/mL) | |
| | 1,2,3,4-Tetrachlorodibenzo- <i>p</i> -dioxin ($^{13}\text{C}_6$, 99%) | 2.5 | 500 |
| | 2,2',3,3',4,5,5',6,6'-Nonachlorobiphenyl ($^{13}\text{C}_{12}$, 99%) | 208 | 2000 |
| | 3,3',4,4'-Tetrabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 77 | 1500 |
| | 2,2',3,4,4',6-Hexabromodiphenyl ether ($^{13}\text{C}_{12}$, 99%) | 139 | 1500 |
| NEW ES-5496 Phenolic Native PAR Standard | | | |
| | | | |
| Unlabeled | | (ng/mL) | |
| | 4-Methoxy-2,3,3',4',5-pentachlorobiphenyl | 1 | |
| | 4-Methoxy-2,2',3,4',5,5'-hexachlorobiphenyl | 1 | |
| | 4-Methoxy-2,2',3,4',5,5',6-heptachlorobiphenyl | 1 | |
| | Pentachloroanisole | 1 | |
| | 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) | 1 | |
| | Pentabromoanisole | 1 | |

Human Exposure Analysis Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------------|---|-----------------|
| NEW ES-5535 | JECS Labeled Mixture Solution | 1.2 mL in water |
| <hr/> | | |
| | Labeled | (ng/mL) |
| | 8-Hydroxy-2'-deoxyguanosine (¹⁵ N ₅ , 98%) | 7500 |
| | DL-Cotinine (2',3',4'- ¹³ C ₃ , 99%) | 150 |

| | | |
|--------------------|------------------------------|-----------------|
| NEW ES-5536 | JECS Native Mixture Solution | 1.2 mL in water |
| <hr/> | | |
| | Unlabeled | (ng/mL) |
| | 8-Hydroxy-2'-deoxyguanosine | 500 |
| | Cotinine | 500 |

Pesticides and Chemical Weapon Standards

Essential to agricultural productivity and general pest control, many pesticides have been found to be toxic to humans and animals, and as such have been banned from use in numerous countries. The Stockholm Convention has focused worldwide attention on analysis of the most toxic pesticides. CIL promotes the development of new isotope-labeled legacy and “new use” pesticides to support laboratories using IDMS for the most accurate analytical results.



Pesticide Standards

CIL continues to add to its already extensive inventory of isotopically labeled standards for pesticide and pesticide metabolite analysis. As a result of this development over the past few years, CIL can now present its standards by category, including organochlorine, organophosphorous, carbamate, triazine, or pyrethroid pesticide standards. You can still find the complete listing if you wish to scan through the comprehensive array of standards.

Chlorinated Cyclodiene Pesticide Standards

Chlorinated cyclodiene pesticides account for seven of the compounds governed by the Stockholm Convention. While production and use of these compounds is stringently regulated if not banned outright, their widespread use for decades and persistence in the environment ensures their presence in the environment and biota for years to come. CIL offers a comprehensive selection of the individual standards, as well as a growing list of convenient mixes.

Organochlorine Pesticide Standards

Organochlorinated pesticides, like chlorinated cyclodiene pesticides, are heavily represented in the list of compounds governed by the Stockholm Convention. Also, like chlorinated cyclodiene pesticides, their widespread use for decades and persistence in the environment ensures their presence in the environment and biota for years to come.

Organophosphate (OP) and Carbamate Insecticides

Organophosphates are a large class of contact insecticides that target the insect's nervous system by interfering with the enzyme acetylcholinesterase, disrupting nerve impulses and killing or disabling the insect. Organophosphate insecticides and chemical warfare nerve agents (such as sarin, tabun, soman, and VX) work in the same way, and metabolites of both groups are quite similar. Organophosphates have a cumulative toxic effect to wildlife, so multiple exposures to the chemicals amplifies the toxicity. Carbamates feature the carbamate ester functional group and kill insects by reversibly inactivating the enzyme acetylcholinesterase, similarly to organophosphate pesticides. They are, however, much less stable in the environment and break down rapidly.

Pyrethroid Insecticides

Pyrethroids are synthetic pesticides developed to mimic the effect of naturally occurring pyrethrins. In general, pyrethroids are low in toxicity to mammals and birds, however, they are potentially highly toxic to fish, have high arthropod toxicity, are fast acting, dissolve poorly in water, and break down quickly, especially in direct sunlight. Pyrethroids became popular as consumer insecticides in the 1990s as replacements for older pesticides, like diazinon and Dursban®, which were phased out for environmental and human-health reasons.

Neonicotinoid Pesticides

Neonicotinoids have received a lot of attention in the mainstream press as a possible cause of colony collapse disorder (CCD) in honeybees, which are critical pollinators in many agricultural environments where their viability enables billions of dollars of commerce each year. While CCD may also have other causes, analytical chemists have focused on neonicotinoid pesticides as a potential contributing factor. CIL has produced several neonicotinoid standards, including labeled standards with chloropyridylmethyl, chlorothiazolylmethyl, and tetrahydrofuranylmethyl substituents, as well as major metabolites of those compounds.

Toxaphene Standards

CIL has put considerable effort into developing the first set of ¹³C-labeled toxaphene standards. The list of labeled and unlabeled standards (with Parlar congener #) continues to grow, so visit CIL's website and watch for future product announcements for more details. The new POPs toxaphene mixtures are ideal for researchers interested in primary investigations of the most prevalent congeners.

Triazine Herbicide and Metabolite Standards

Atrazine is one of the most widely used herbicides in the world. In recent years, studies on the correlation of physical and reproductive disorders in frogs with atrazine exposure has been a controversial topic. With CIL's comprehensive collection of carefully purified and prepared standards of atrazine and its many metabolites, researchers should have some powerful tools to refine their investigations.

Pesticide Standard Mixtures

New applications and increased testing by IDMS have led to the development of several pesticides mixtures being offered for the first time in this catalog. Our expanded POPs pesticide calibration series and related spiking mixtures contain all pesticides listed as Stockholm Convention POPs, including kepone (aka chlordcone), HCHs (including lindane), pentachlorobenzene, and endosulfan I and II. These solutions allow analysts to use preformulated mixtures for detection and quantification of the complete series of these important POPs.

Chemical Weapon Metabolite Standards

Often quite similar to metabolites of common pesticides, chemical weapons metabolite standards help researchers determine potential contamination from dangerous compounds, such as nerve agents and other toxic chemicals. Several metabolites, degradation byproducts, and others are represented in this section.

Chlorinated Cyclodiene Pesticide Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-----------------|--|---|---|----------------------|
| CLM-4725-1.2 | Aldrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7441-1.2 | Aldrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-8087-1.2 | cis-Chlordane (α) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2419-1.2 | cis-Chlordane (α) (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane |
| | ULM-2419-25 | | neat | 25 mg |
| CLM-4792-1.2 | trans-Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2420-1.2 | trans-Chlordane (γ) (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2420-25 | | neat | 25 mg |
| | CLM-4814-1.2 | Chlordecone (kepone) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane |
| ULM-2301-1.2 | Chlordecone (kepone) (unlabeled) | $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2301-0.1 | | neat | 0.1 g |
| | CLM-4758-1.2 | Chlordene ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane |
| ULM-7443-1.2 | Chlordene (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4726-1.2 | Dieldrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7230-1.2 | Dieldrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-6025-1.2 | Endosulfan I ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| DLM-2862-1.2 | Endosulfan I (D_4 , 97%) | $\text{C}_9\text{D}_4\text{H}_2\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7447-1.2 | Endosulfan I (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-6026-1.2 | Endosulfan II ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7448-1.2 | Endosulfan II (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7531-1.2 | Endosulfan sulfate ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_4\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7990-1.2 | Endosulfan sulfate (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_4\text{S}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4782-1.2 | Endrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7444-1.2 | Endrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | CLM-4815-1.2 | Endrin aldehyde ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_{10}\text{Cl}_6\text{O}$ | 100 µg/mL in nonane |
| | CLM-4815-50 | | neat | 50 µg |
| NEW | ULM-8958-1.2 | Endrin aldehyde (unlabeled) | $\text{C}_{12}\text{H}_{10}\text{Cl}_6\text{O}$ | 100 µg/mL in nonane |
| | ULM-8958-50 | | neat | 50 µg |
| NEW | CLM-4816-1.2 | Endrin ketone ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane |
| | CLM-4816-50 | | neat | 50 µg |
| NEW | ULM-8956-1.2 | Endrin ketone (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane |
| | ULM-8956-50 | | neat | 50 µg |
| CLM-4759-1.2 | Heptachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_7$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-1.2 | Heptachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2424-0.1 | | neat | 0.1 g |
| | CLM-4734-1.2 | cis-Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane |
| ULM-2425-1.2 | cis-Heptachlor epoxide (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2425-0.1 | | neat | 0.1 g |
| ULM-7869-1.2 | trans-Heptachlor epoxide (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4727-1.2 | Isodrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7442-1.2 | Isodrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4814-1.2 | Kepone (chlordecone) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-1.2 | Kepone (chlordecone) (unlabeled) | $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW | ULM-2301-0.1 | | neat | 0.1 g |
| | CLM-4813-1.2 | Mirex ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{12}$ | 100 µg/mL in nonane |
| NEW | CLM-2078-1 | Mirex ($^{13}\text{C}_8$, 99%) | * $\text{C}_8\text{C}_2\text{Cl}_{12}$ | 200 µg/mL in toluene |
| | ULM-2427-1.2 | Mirex (unlabeled) | $\text{C}_{10}\text{Cl}_{12}$ | 100 µg/mL in nonane |
| CLM-4811-1.2 | cis-Nonachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7445-1.2 | cis-Nonachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4735-1.2 | trans-Nonachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7229-1.2 | trans-Nonachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4729-1.2 | Oxychlordane ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_4\text{Cl}_8\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-1.2 | Oxychlordane (unlabeled) | $\text{C}_{10}\text{H}_4\text{Cl}_8\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-SM-1.2 | | | 100 µg/mL in methanol | 1.2 mL |

NOTE: Some standards also available in less than uniformly labeled forms. Please inquire if interested.

Organochlorine (OC) Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount | |
|--------------|---|---|--------------------------------------|---------------------|--------|
| CLM-4725-1.2 | Aldrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7441-1.2 | Aldrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-2482-1.2 | α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7232-1.2 | α -HCH (α -BHC) (unlabeled) | $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-3623-1.2 | β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_6\text{Cl}_6$ | 50 µg/mL in nonane | 2 × 1.2 mL | |
| ULM-6132-1.2 | β -HCH (β -BHC) (unlabeled) | $\text{C}_6\text{H}_6\text{Cl}_6$ | 50 µg/mL in nonane | 2 × 1.2 mL | |
| CDLM-624-1.2 | γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%; D_6 , 99%) | * $\text{C}_6\text{D}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1282-1.2 | γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-6133-1.2 | γ -HCH (γ -BHC) (lindane) (unlabeled) | $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-3648-1.2 | δ -HCH (δ -BHC) ($^{13}\text{C}_6$, 99%) | * $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7233-1.2 | δ -HCH (δ -BHC) (unlabeled) | $\text{C}_6\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-8087-1.2 | cis-Chlordane (α) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL | |
| NEW | ULM-2419-1.2 | cis-Chlordane (α) (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-2419-25 | | neat | 25 mg | |
| CLM-4792-1.2 | trans-Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL | |
| NEW | ULM-2420-1.2 | trans-Chlordane (γ) (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_8$ | 100 µg/mL in nonane | 1.2 mL |
| | ULM-2420-25 | | neat | 25 mg | |
| CLM-4814-1.2 | Chlordecone (kepone) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-2301-1.2 | Chlordecone (kepone) (unlabeled) | $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-2301-0.1 | | neat | 0.1 g | | |
| CLM-4758-1.2 | Chlordene ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7443-1.2 | Chlordene (unlabeled) | $\text{C}_{10}\text{H}_6\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-6999-1.2 | 2,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) [(o,p'-Dichlorodiphenyl) dichloroethane] | * $\text{C}_{12}\text{C}_2\text{H}_{10}\text{Cl}_4$ | 50 µg/mL in nonane | 1.2 mL | |
| ULM-7450-1.2 | 2,4'-DDD (unlabeled) [(o,p'-dichlorodiphenyl) dichloroethane] | $\text{C}_{14}\text{H}_{10}\text{Cl}_4$ | 50 µg/mL in nonane | 1.2 mL | |
| CLM-7100-1.2 | 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) [(p,p'-dichlorodiphenyl) dichloroethane] | * $\text{C}_{12}\text{C}_2\text{H}_{10}\text{Cl}_4$ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-3533-1.2 | 4,4'-DDD (ring-D ₈ , 98%) [(p,p'-dichlorodiphenyl) dichloroethane] | $\text{C}_{14}\text{D}_8\text{H}_2\text{Cl}_4$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7216-1.2 | 4,4'-DDD (unlabeled) [(p,p'-dichlorodiphenyl) dichloroethane] | $\text{C}_{14}\text{H}_{10}\text{Cl}_4$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-4693-1.2 | 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) [(o,p'-dichlorodiphenyl) dichloroethylene] | (Cl* $\text{C}_6\text{H}_4)_2\text{C}=\text{CCl}_2$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-6251-1.2 | 2,4'-DDE (unlabeled) [(o,p'-dichlorodiphenyl) dichloroethylene] | $\text{C}_{14}\text{H}_8\text{Cl}_4$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1627-1.2 | 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) [(p,p'-dichlorodiphenyl) dichloroethylene] | (Cl* $\text{C}_6\text{H}_4)_2\text{C}=\text{CCl}_2$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1627-5 | | neat | 5 mg | | |
| ULM-6137-1.2 | 4,4'-DDE (unlabeled) [(p,p'-dichlorodiphenyl) dichloroethylene] | (Cl $\text{C}_6\text{H}_4)_2\text{C}=\text{CCl}_2$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-4692-1.2 | 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) [(o,p'-dichlorodiphenyl) trichloroethane] | (Cl* $\text{C}_6\text{H}_4)_2\text{CHCCl}_3$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-6134-1.2 | 2,4'-DDT (unlabeled) [(o,p'-dichlorodiphenyl) trichloroethane] | Cl $\text{C}_6\text{H}_4\text{CH}(\text{CCl}_3)\text{C}_6\text{H}_4\text{Cl}$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1281-1.2 | 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) [(p,p'-dichlorodiphenyl) trichloroethane] | (Cl* $\text{C}_6\text{H}_4)_2\text{CHCCl}_3$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1281-5 | | neat | 5 mg | | |
| ULM-6135-1.2 | 4,4'-DDT (unlabeled) [(p,p'-dichlorodiphenyl) trichloroethane] | (Cl $\text{C}_6\text{H}_4)_2\text{CHCCl}_3$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-816-1.2 | 2,6-Dichloro-4-nitroaniline (ring- $^{13}\text{C}_6$, 99%) | $\text{Cl}_2^*\text{C}_6\text{H}_2(\text{NO}_2)\text{NH}_2$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-4726-1.2 | Dieldrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7230-1.2 | Dieldrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-6025-1.2 | Endosulfan I ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| DLM-2862-1.2 | Endosulfan I (D ₄ , 97%) | $\text{C}_9\text{D}_4\text{H}_2\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7447-1.2 | Endosulfan I (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-6026-1.2 | Endosulfan II ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7448-1.2 | Endosulfan II (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_3\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-7531-1.2 | Endosulfan sulfate ($^{13}\text{C}_9$, 99%) | * $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_4\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7990-1.2 | Endosulfan sulfate (unlabeled) | $\text{C}_9\text{H}_6\text{Cl}_6\text{O}_4\text{S}$ | 100 µg/mL in nonane | 1.2 mL | |

Organochlorine (OC) Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|-----------------------|--------|
| CLM-4782-1.2 | Endrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7444-1.2 | Endrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4815-1.2 | Endrin aldehyde ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_{10}\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4815-50 | | | neat | 50 µg |
| NEW ULM-8958-1.2 | Endrin aldehyde (unlabeled) | $\text{C}_{12}\text{H}_{10}\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8958-50 | | | neat | 50 µg |
| NEW CLM-4816-1.2 | Endrin ketone ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4816-50 | | | neat | 50 µg |
| NEW ULM-8956-1.2 | Endrin ketone (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8956-50 | | | neat | 50 µg |
| CLM-4759-1.2 | Heptachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_7$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-1.2 | Heptachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-0.1 | | | neat | 0.1 g |
| CLM-4734-1.2 | <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2425-1.2 | <i>cis</i> -Heptachlor epoxide (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2425-0.1 | | | neat | 0.1 g |
| ULM-7869-1.2 | <i>trans</i> -Heptachlor epoxide (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_7\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-351-1.2 | Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | * C_6Cl_6 | 100 µg/mL in nonane | 1.2 mL |
| ULM-6130-1.2 | Hexachlorobenzene (unlabeled) | C_6Cl_6 | 100 µg/mL in nonane | 1.2 mL |
| CLM-4727-1.2 | Isodrin ($^{13}\text{C}_{12}$, 99%) | * $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7442-1.2 | Isodrin (unlabeled) | $\text{C}_{12}\text{H}_8\text{Cl}_6$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4814-1.2 | Kepone (chlordecone) ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-1.2 | Kepone (chlordecone) (unlabeled) | $\text{C}_{10}\text{Cl}_{10}\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-0.1 | | | neat | 0.1 g |
| CLM-4683-1.2 | Methoxychlor (ring- $^{13}\text{C}_{12}$, 99%) | ($\text{CH}_3\text{O}^*\text{C}_6\text{H}_4)_2\text{CHCCl}_3$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7440-1.2 | Methoxychlor (unlabeled) | ($\text{CH}_3\text{OC}_6\text{H}_4)_2\text{CHCCl}_3$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4813-1.2 | Mirex ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{Cl}_{12}$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-2078-1 | Mirex ($^{13}\text{C}_8$, 99%) | * $\text{C}_8\text{C}_2\text{Cl}_{12}$ | 100 µg/mL in toluene | 1 mL |
| ULM-2427-1.2 | Mirex (unlabeled) | $\text{C}_{10}\text{Cl}_{12}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2427-0.1 | | | neat | 0.1 g |
| CLM-4811-1.2 | <i>cis</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7445-1.2 | <i>cis</i> -Nonachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4735-1.2 | <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7229-1.2 | <i>trans</i> -Nonachlor (unlabeled) | $\text{C}_{10}\text{H}_5\text{Cl}_9$ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4729-1.2 | Oxychlordane ($^{13}\text{C}_{10}$, 99%) | * $\text{C}_{10}\text{H}_4\text{Cl}_8\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-1.2 | Oxychlordane (unlabeled) | $\text{C}_{10}\text{H}_4\text{Cl}_8\text{O}$ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-SM-1.2 | | | 100 µg/mL in methanol | 1.2 mL |

Organophosphate (OP) Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------|--|------------------------------------|---|---------|
| DLM-6000-1.2 | Acephate (D_6 , 98%) | $C_4D_6H_4NO_3PS$ | 100 $\mu\text{g/mL}$ in acetonitrile- D_3 | 1.2 mL |
| ULM-7263-1.2 | Acephate (unlabeled) | $C_4H_{10}NO_3PS$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| CDNLM-6786-1.2 | Aminomethylphosphonic acid (AMPA) | $*CH_4D_2^*NO_3P$ | 100 $\mu\text{g/mL}$ in H_2O | 1.2 mL |
| NEW CDNLM-6786-10 | (^{13}C , 99%; ^{15}N , 98%, methylene- D_2 , 98%) | | 100 $\mu\text{g/mL}$ in H_2O | 10 mL |
| NEW ULM-9399-1.2 | Azinphos-methyl (unlabeled) | $C_{10}PN_3H_{12}S_2O_3$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| DLM-7152 | Bensulide (isopropoxy- D_{14} , 98%) | $C_{14}D_{14}H_{10}NO_4PS_3$ | | Inquire |
| DLM-4360-1.2 | Chlorpyrifos (diethyl- D_{10} , 99%) | $C_9D_{10}HCl_3NO_3PS$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-7489-1.2 | Chlorpyrifos (unlabeled) | $C_9H_{11}Cl_3NO_3PS$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW DLM-7153-1.2 | Chlorpyrifos-methyl (dimethyl- D_6 , 98%) | $C_7HCl_3D_6NO_3PS$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW ULM-9538-1.2 | Chlorpyrifos-methyl (unlabeled) | $C_7H_7Cl_3NO_3PS$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| DLM-1148-1.2 | Diazinon (diethyl- D_{10} , 98%) | $C_{12}H_{11}D_{10}N_2O_3PS$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW DLM-1148-5 | | | neat | 5 mg |
| NEW DLM-1148-A-1.2 | | | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-6575-A-1.2 | Diazinon (unlabeled) | $C_{12}H_{21}N_2O_3PS$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-6575-S-10X-1.2 | | | 1000 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| DLM-2829-0.01 | Dichlorvos (dimethyl- D_6 , 98%) | $C_4D_6HCl_2O_4P$ | neat | 10 mg |
| ULM-7217-1.2 | Dichlorvos (unlabeled) | $(H_3CO)_2POOCH=CCl_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW DLM-4851-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) (diethyl- D_{10} , 98%) | $C_4D_{10}KO_4P$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9287-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) (unlabeled) | $C_4H_{10}KO_4P$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-4852-1.2 | O,O-Diethyl thiophosphate, potassium salt (DETP) (diethyl- D_{10} , 98%) | $C_4D_{10}KO_3PS$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ERD-119 | O,O-Diethyl thiophosphate, potassium salt (DETP) (unlabeled) | $C_4H_{10}KO_3PS$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW DLM-9003-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) (diethyl- D_{10} , 98%) | $C_4D_{10}KO_2PS_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9002-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) (unlabeled) | $C_4H_{10}KO_2PS_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ULM-9898-1.2 | Diisopropyl methylphosphonate (unlabeled) | $C_7H_{17}O_3P$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-7151-1.2 | Dimethoate (O,O-dimethyl- D_6 , 98%) | $C_5D_6H_6NO_3PS_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-7972-1.2 | Dimethoate (unlabeled) | $C_5H_{12}NO_3PS_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| NEW DLM-8868-1.2 | O,O-Dimethyl phosphate, potassium salt (DMP) (dimethyl- D_6 , 98%) | $C_2D_6KO_4P$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-8867-1.2 | O,O-Dimethyl phosphate, potassium salt (DMP) (unlabeled) | $C_2H_6KO_3PS$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW DLM-8904-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) (dimethyl- D_6 , 98%) CP 97% | $C_2D_6KO_3PS$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-8905-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) (unlabeled) CP 97% | $C_2H_6KO_3PS$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW DLM-4541-M-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) (dimethyl- D_6 , 98%) | $C_2D_6KO_2PS_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9004-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) (unlabeled) | $C_2H_6KO_2PS_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ULM-9899-1.2 | Dipinacolyl methylphosphonate (unlabeled) | $C_7H_{17}O_3P$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-7183 | Disulfoton (O,O-diethyl- D_{10} , 98%) | $C_8D_{10}H_9O_2PS_3$ | | Inquire |
| ULM-6091-1.2 | Ethyl dimethylamidophosphate, sodium salt (unlabeled) | $C_4H_{11}NO_3PNa$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-6098-1.2 | Ethyl hydrogen methylphosphonate (ethyl- D_5 , 98%) | $C_3H_4D_5O_3P$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ULM-6099-1.2 | Ethyl methylphosphonic acid (unlabeled) | $C_3H_9O_3P$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-2878-0.01 | Fenitrothion (O,O-dimethyl- D_6 , 98%) | $C_9D_6H_6NO_5PS$ | neat | 10 mg |
| CLM-4545-1.2 | Fonofos (ring- $^{13}\text{C}_6$, 99%) | $*C_6C_4H_{15}OPS_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-6694-1.2 | Fonofos (unlabeled) | $CH_2CH_3P(S)(OCH_2CH_3)(SC_6H_5)$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |

Organophosphate (OP) Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------|--|---|--------------------------------------|--------|
| CNLM-4666-1.2 | Glyphosate ($2\text{-}^{13}\text{C}$, 99%; ^{15}N , 98%+) CP 96% | $\text{HOO}^*\text{CCH}_2\text{NHCH}_2\text{PO(OH)}_2$ | 100 $\mu\text{g/mL}$ in water | 1.2 mL |
| CNLM-4666-10 | | | 100 $\mu\text{g/mL}$ in water | 10 mL |
| CNLM-4666-10X-1.2 | | | 1000 $\mu\text{g/mL}$ in water | 1.2 mL |
| ULM-6876-1.2 | Glyphosate (unlabeled) | $\text{HOOCCH}_2\text{NHCH}_2\text{PO(OH)}_2$ | 100 $\mu\text{g/mL}$ in water | 1.2 mL |
| ULM-6093-1.2 | Isopropyl methylphosphonic acid (unlabeled) | $\text{C}_4\text{H}_{11}\text{O}_3\text{P}$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-4476-1.2 | Malathion (D_{10} , 99%) | $\text{C}_{10}\text{D}_{10}\text{H}_9\text{O}_6\text{PS}_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-8122-1.2 | Malathion (unlabeled) | $\text{C}_{10}\text{H}_{15}\text{O}_6\text{PS}_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| NEW CLM-9050-1.2 | Malathion diacid ($^{13}\text{C}_4$, 99%) CP 97% | $^*\text{C}_6\text{H}_{11}\text{O}_6\text{PS}_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9073-1.2 | Malathion diacid (unlabeled) | $\text{C}_4\text{C}_2\text{H}_{11}\text{O}_6\text{PS}_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW DLM-7149-1.2 | Methamidophos (dimethyl- D_6 , 98%) | $\text{C}_2\text{D}_6\text{H}_2\text{NO}_2\text{PS}$ | 100 $\mu\text{g/mL}$ in dioxane | 1.2 mL |
| NEW ULM-8872-1.2 | Methamidophos (unlabeled) | $\text{C}_2\text{H}_8\text{NO}_2\text{PS}$ | 100 $\mu\text{g/mL}$ in dioxane | 1.2 mL |
| CDLM-6100-1.2 | Methylphosphonic acid (^{13}C , 99%; methyl- D_3 , 98%) | $^*\text{CD}_3\text{H}_2\text{O}_3\text{P}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-6196-1.2 | Methylphosphonic acid (methyl- D_3 , 98%) | $\text{CD}_3\text{H}_2\text{O}_3\text{P}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| ERM-038 | Methylphosphonic acid (unlabeled) | $\text{CH}_3\text{P(O)(OH)}_2$ | 1000 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| CLM-6620-1.2 | Methylphosphonic acid, mono-(1,2,2-trimethylpropyl) ester (trimethylpropyl- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{CH}_{17}\text{O}_3\text{P}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| DLM-7150-1.2 | Oxydemeton methyl (O,O -dimethyl- D_6 , 98%) | $\text{C}_6\text{D}_6\text{H}_9\text{O}_4\text{PS}_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-8579-1.2 | Oxydemeton methyl (unlabeled) | $\text{C}_6\text{H}_{15}\text{O}_4\text{PS}_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| DLM-2970-1.2 | Parathion (diethyl- D_{10} , 98%) | $\text{NO}_2(\text{C}_6\text{H}_4)\text{OP(=S)(OC}_2\text{D}_5)_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| ULM-8144-1.2 | Parathion (unlabeled) | $\text{NO}_2(\text{C}_6\text{H}_4)\text{OP(=S)(OC}_2\text{H}_5)_2$ | 100 $\mu\text{g/mL}$ in nonane | 1.2 mL |
| CLM-4544-1.2 | Phorate (diethoxy- $^{13}\text{C}_4$, 99%) | $(^*\text{C}_2\text{H}_5\text{O})_2\text{P(S)SC}_2\text{SC}_2\text{H}_5$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-7567-1.2 | Phorate (unlabeled) | $(\text{C}_2\text{H}_5\text{O})_2\text{P(S)SC}_2\text{SC}_2\text{H}_5$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| DLM-4667-1.2 | Phosmet (dimethyl- D_6 , 98%) | $\text{C}_{11}\text{D}_6\text{H}_9\text{NO}_4\text{PS}_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| ULM-8454-1.2 | Phosmet (unlabeled) | $\text{C}_{11}\text{H}_{12}\text{NO}_4\text{PS}_2$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| CLM-4543 | Terbufos (diethoxy- $^{13}\text{C}_4$, 99%) | $\text{C}(\text{CH}_3)_3\text{SCH}_2\text{SP(S)(O}^*\text{CH}_2\text{*CH}_3)_2$ | Inquire | |
| NEW CLM-9049-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (4,5,6- $^{13}\text{C}_3$, 99%) CP 97% | $^*\text{C}_3\text{C}_2\text{H}_2\text{Cl}_3\text{NO}$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |
| NEW ULM-9204-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (unlabeled) | $\text{C}_5\text{H}_2\text{Cl}_3\text{NO}$ | 100 $\mu\text{g/mL}$ in acetonitrile | 1.2 mL |

Neonicotinoid Insecticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------|--|---|----------------------------------|--------|
| NEW CLM-9653-1.2 | Acetamiprid (pyridylmethyl- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_4\text{H}_{11}\text{ClN}_4$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9734-1.2 | Acetamiprid (unlabeled) | $\text{C}_{10}\text{H}_{11}\text{ClN}_4$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CLM-9598-1.2 | 6-Chloronicotinic acid ($^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{H}_4\text{ClNO}_2$ | 100 $\mu\text{g/mL}$ in MTBE | 1.2 mL |
| NEW ULM-9604-1.2 | 6-Chloronicotinic acid (unlabeled) | $\text{C}_6\text{H}_4\text{ClNO}_2$ | 100 $\mu\text{g/mL}$ in MTBE | 1.2 mL |
| NEW CNLM-9940-1.2 | Clothianidin (thiazole- $^{13}\text{C}_3$, 99%; ^{15}N , 98%) | $^*\text{C}_3\text{C}_3\text{H}_8\text{Cl}^*\text{NN}_4\text{O}_2\text{S}$ | Inquire | |
| NEW ULM-9941-1.2 | Clothianidin (unlabeled) | $\text{C}_6\text{H}_8\text{ClN}_5\text{O}_2\text{S}$ | Inquire | |
| NEW CLM-9594-1.2 | Dinotefuran (furylmethyl- $^{13}\text{C}_5$, 99%) | $^*\text{C}_5\text{C}_2\text{H}_{14}\text{N}_4\text{O}_3$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9732-1.2 | Dinotefuran (unlabeled) | $\text{C}_7\text{H}_{14}\text{N}_4\text{O}_3$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW DLM-8512-1.2 | Imidacloprid (4,4,5,5- D_4 , 98%) | $\text{C}_9\text{H}_6\text{D}_4\text{ClN}_5\text{O}_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-8513-1.2 | Imidacloprid (unlabeled) | $\text{C}_9\text{H}_{10}\text{ClN}_5\text{O}_2$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CNLM-9869-1.2 | Sulfoxaflor (cyano- ^{13}C , 99%; cyano- ^{15}N , 98%) | $^*\text{CC}_9\text{H}_{10}\text{F}_3^*\text{N}_2\text{NOS}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9870-1.2 | Sulfoxaflor (unlabeled) | $\text{C}_{10}\text{H}_{10}\text{F}_3\text{N}_2\text{NOS}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CLM-9690-1.2 | 3-Tetrahydrofuroic acid ($^{13}\text{C}_5$, 99%) | $^*\text{C}_5\text{H}_8\text{O}_3$ | 100 $\mu\text{g/mL}$ in MTBE | 1.2 mL |
| NEW ULM-9691-1.2 | 3-Tetrahydrofuroic acid (unlabeled) | $\text{C}_5\text{H}_8\text{O}_3$ | 100 $\mu\text{g/mL}$ in MTBE | 1.2 mL |
| NEW CLM-9652-1.2 | Thiacloprid (pyridylmethyl- $^{13}\text{C}_6$, 99%) | $^*\text{C}_6\text{C}_4\text{H}_9\text{ClN}_4\text{S}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW ULM-9733-1.2 | Thiacloprid (unlabeled) | $\text{C}_{10}\text{H}_9\text{ClN}_4\text{S}$ | 100 $\mu\text{g/mL}$ in methanol | 1.2 mL |
| NEW CNLM-9860-1.2 | Thiamethoxam (thiazole- $^{13}\text{C}_3$, 99%; ^{15}N , 98%) | $^*\text{C}_3\text{C}_5\text{H}_{10}\text{Cl}^*\text{NN}_4\text{O}_3\text{S}$ | Inquire | |
| NEW ULM-9939-1.2 | Thiamethoxam (unlabeled) | $\text{C}_8\text{H}_{10}\text{ClN}_5\text{O}_3\text{S}$ | Inquire | |

Carbamate Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------------|---|---|--------------------------------|---------|
| NEW CDLM-9820-1.2 | Aldicarb (¹³ C ₂ , 98%; D ₃ , 98%) | C ₅ *C ₂ H ₁₁ D ₃ N ₂ O ₂ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9823-1.2 | Aldicarb (unlabeled) | C ₇ H ₁₄ N ₂ O ₂ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7140 | Bendiocarb (¹³ C ₃ , 99%) | *C ₃ C ₈ H ₁₃ NO ₄ | | Inquire |
| ULM-8638 | Bendiocarb (unlabeled) | C ₁₁ H ₁₃ NO ₄ | | Inquire |
| CLM-4682-1.2 | Carbaryl (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₁ NO ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-8096-1.2 | Carbaryl (unlabeled) | C ₁₀ H ₇ CO ₂ NHCH ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1911-1.2 | Carbofuran (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₅ NO ₃ | 100 µg/mL in <i>p</i> -dioxane | 1.2 mL |
| ULM-7419-1.2 | Carbofuran (unlabeled) | C ₁₂ H ₁₅ NO ₃ | 100 µg/mL in <i>p</i> -dioxane | 1.2 mL |
| ULM-6875-1.2 | Carbofuran phenol (unlabeled) | C ₁₀ H ₁₂ O ₂ | 200 µg/mL in nonane | 1.2 mL |
| CNLM-7148-1.2 | Methomyl (acetohydroxamate- ¹³ C ₂ , 99%; ¹⁵ N, 98%) | *C ₂ C ₃ H ₁₀ N*NO ₂ S | 100 µg/mL in methanol | 1.2 mL |
| ULM-8639-1.2 | Methomyl (unlabeled) | C ₅ H ₁₀ NNO ₂ S | 100 µg/mL in methanol | 1.2 mL |
| DLM-7141-1.2 | Propoxur (isopropyl-D ₇ , 98%) | C ₁₁ D ₇ H ₈ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9765-1.2 | Propoxur (unlabeled) | C ₁₁ H ₁₅ NO ₃ | 100 µg/mL in nonane | 1.2 mL |

Triazine Herbicide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------------|--|---|---|---------|
| CLM-8316-1.2 | Ammeline (desethyldeisopropylhydroxyatrazine) (ring- ¹³ C ₃ , 99%) | NH ₂ (*C ₃ N ₃ OH)NH ₂ | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8323-1.2 | Ammeline (desethyldeisopropylhydroxyatrazine) (unlabeled) | NH ₂ (C ₃ N ₃ OH)NH ₂ | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| CLM-3737-1.2 | Atrazine (ring- ¹³ C ₃ , 99%) | (CH ₃) ₂ CHNH(*C ₃ N ₃ Cl)NHCH ₂ CH ₃ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1149-1.2 | Atrazine (ethylamine-D ₅ , 98%) | (CH ₃) ₂ CHNH(C ₃ N ₃ Cl)NHCD ₂ CD ₃ | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-1149-5 | | | 100 µg/mL in nonane | 5 mg |
| ULM-7235-1.2 | Atrazine (unlabeled) | (CH ₃) ₂ CHNH(C ₃ N ₃ Cl)NHCH ₂ CH ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3894-1.2 | Atrazine mercapturate (ring- ¹³ C ₃ , 99%) | *C ₃ C ₁₀ H ₂₂ N ₆ O ₃ S | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7346-1.2 | Atrazine mercapturate (unlabeled) | C ₁₃ H ₂₂ N ₆ O ₃ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8311-1.2 | Atrazinethiol (ring- ¹³ C ₃ , 99%) | (CH ₃ CH ₂ NH)*C ₃ N ₃ (SH)(NHCH(CH ₃) ₂) | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8318-1.2 | Atrazinethiol (unlabeled) | (CH ₃ CH ₂ NH)C ₃ N ₃ (SH)(NHCH(CH ₃) ₂) | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8313-1.2 | Desethylatrazine (ring- ¹³ C ₃ , 99%) CP 97% | (CH ₃) ₂ CHNH(*C ₃ N ₃ Cl)NH ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8320-1.2 | Desethylatrazine (unlabeled) | (CH ₃) ₂ CHNH(C ₃ N ₃ Cl)NH ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7528-1.2 | Desethyl desisopropyl atrazine (¹³ C ₃ , 99%) CP 95% | *C ₃ H ₄ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8001-1.2 | Desethyl desisopropyl atrazine (unlabeled) | C ₃ H ₄ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8315-1.2 | Desethylhydroxyatrazine (ring- ¹³ C ₃ , 99%) | (CH ₂ N)*C ₃ N ₃ (OH)(NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8322-1.2 | Desethylhydroxyatrazine (unlabeled) | (CH ₂ N)C ₃ N ₃ (OH)(NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| CLM-8312-1.2 | Desisopropylatrazine (ring- ¹³ C ₃ , 99%) | CH ₃ CH ₂ NH(*C ₃ N ₃ Cl)NH ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8319-1.2 | Desisopropylatrazine (unlabeled) | CH ₃ CH ₂ NH(C ₃ N ₃ Cl)NH ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8314-1.2 | Desisopropylhydroxyatrazine (ring- ¹³ C ₃ , 99%) | (CH ₃ CH ₂ NH)*C ₃ N ₃ (OH)(NH ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8321-1.2 | Desisopropylhydroxyatrazine (unlabeled) | (CH ₃ CH ₂ NH)C ₃ N ₃ (OH)(NH ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| CLM-8310-1.2 | Hydroxyatrazine (ring- ¹³ C ₃ , 99%) | (CH ₃ CH ₂ NH)*C ₃ N ₃ (OH)(NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8317-1.2 | Hydroxyatrazine (unlabeled) | (CH ₃ CH ₂ NH)C ₃ N ₃ (OH)(NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| CLM-3738-1.2 | Propazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₆ H ₁₆ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8304 | Propazine (unlabeled) | C ₉ H ₁₆ CIN ₅ | | Inquire |
| CLM-3739-1.2 | Simazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₄ H ₁₂ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-3739-A-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7893-1.2 | Simazine (unlabeled) | C ₇ H ₁₂ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-7893-A-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |

Pyrethroid Pesticide and Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------------|---|--|--|--------|
| CLM-7293-1.2 | Cyfluthrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7454-1.2 | Cyfluthrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7292-1.2 | Cypermethrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7453-1.2 | Cypermethrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| NEW CDLM-9205-1.2 | cis-DCCA (1,carboxyl- ¹³ C ₂ , 99%; 1-D, 97%) | C ₆ *C ₂ H ₉ Cl ₂ O ₂ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| NEW ULM-9176-1.2 | cis-DCCA (unlabeled) | C ₈ H ₁₀ Cl ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CDLM-9206-1.2 | trans-DCCA (1,carboxyl- ¹³ C ₂ , 99%; 1-D, 97%) CP 97% | C ₆ *C ₂ H ₉ Cl ₂ O ₂ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| NEW ULM-9175-1.2 | trans-DCCA (unlabeled) | C ₈ H ₁₀ Cl ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7389-1.2 | 4-Fluoro-3-phenoxybenzoic acid (¹³ C ₆ , 99%) | *C ₆ C ₇ H ₉ FO ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7391-1.2 | 4-Fluoro-3-phenoxybenzoic acid (unlabeled) | C ₁₃ H ₉ FO ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7322-1.2 | cis-Permethrin (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8526-1.2 | cis-Permethrin (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| CLM-7323-1.2 | trans-Permethrin (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8527-1.2 | trans-Permethrin (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| CLM-4542-1.2 | 3-Phenoxybenzoic acid (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |
| CLM-4542-SA-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-6781-1.2 | 3-Phenoxybenzoic acid (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |
| ULM-6781-SA-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |

Toxaphene Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|--|--------------------|--------|
| NEW ULM-9429-1.2 | Hp-Sed (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9428-1.2 | Hx-Sed (unlabeled) | C ₁₀ H ₁₂ Cl ₆ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7930-1.2 | Parlar 26 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7828-1.2 | Parlar 26 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8705-1.2 | Parlar 32 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8665-1.2 | Parlar 32 (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9005-1.2 | Parlar 38 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8719-1.2 | Parlar 39 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8767-1.2 | Parlar 39 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9430-1.2 | Parlar 40 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9431-1.2 | Parlar 41 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9432-1.2 | Parlar 44 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7931-1.2 | Parlar 50 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7829-1.2 | Parlar 50 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7932-1.2 | Parlar 62 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7830-1.2 | Parlar 62 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8720-1.2 | Parlar 69 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8768-1.2 | Parlar 69 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8721-1.2 | Parlar 70 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8769-1.2 | Parlar 70 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |

See page 226 for mixtures containing toxaphene congeners.

Individual Pesticide and Pesticide Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------------|--|---|--|------------|
| DLM-6000-1.2 | Acephate (D ₆ , 98%) | C ₄ D ₆ H ₄ NO ₃ PS | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| ULM-7263-1.2 | Acephate (unlabeled) | C ₄ H ₁₀ NO ₃ PS | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CLM-9653-1.2 | Acetamiprid (pyridylmethyl- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₁ CIN ₄ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9734-1.2 | Acetamiprid (unlabeled) | C ₁₀ H ₁₁ CIN ₄ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-4546-1.2 | Acetochlor (ring- ¹³ C ₆ , 99%) | *C ₆ C ₈ H ₂₀ CINO ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9824-1.2 | Acetochlor (unlabeled) | C ₆ C ₈ H ₂₀ CINO ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-3727-1.2 | Alachlor (ring- ¹³ C ₆ , 99%) CP 96% | *C ₆ C ₈ H ₂₀ CINO ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3687-1.2 | Alachlor acetyl cysteine adduct (ring- ¹³ C ₆ , 99%) | *C ₆ C ₁₃ H ₂₈ N ₂ O ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CDLM-9820-1.2 | Aldicarb (¹³ C ₂ , 98%; D ₃ , 98%) | C ₅ *C ₂ H ₁₁ D ₃ N ₂ O ₂ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9823-1.2 | Aldicarb (unlabeled) | C ₇ H ₁₄ N ₂ O ₂ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-4725-1.2 | Aldrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7441-1.2 | Aldrin (unlabeled) | C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CDNLM-6786-1.2 | Aminomethylphosphonic acid (AMPA) | *CH ₄ D ₂ *NO ₃ P | 100 µg/mL in H ₂ O | 1.2 mL |
| NEW CDNLM-6786-10 | (¹³ C, 99%; ¹⁵ N, 98%, methylene-D ₂ , 98%) | | | 10 mL |
| CLM-8316-1.2 | Ammeline (desethyl desisopropylhydroxyatrazine) (ring- ¹³ C ₃ , 99%) | NH ₂ (*C ₃ N ₃ OH)NH ₂ | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8323-1.2 | Ammeline (desethyl desisopropylhydroxyatrazine) (unlabeled) | NH ₂ (C ₃ N ₃ OH)NH ₂ | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| CLM-3737-1.2 | Atrazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₅ H ₁₄ CIN ₅ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1149-1.2 | Atrazine (ethylamine-D ₅ , 98%) | C ₈ H ₉ D ₅ CIN ₅ | 100 µg/mL in nonane | 1.2 mL |
| DLM-1149-5 | | | neat | 5 mg |
| ULM-7235-1.2 | Atrazine (unlabeled) | C ₈ H ₁₄ CIN ₅ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3894-1.2 | Atrazine mercapturate (ring- ¹³ C ₃ , 99%) | *C ₃ C ₁₀ H ₂₂ N ₆ O ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7346-1.2 | Atrazine mercapturate (unlabeled) | C ₁₃ H ₂₂ N ₆ O ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8311-1.2 | Atrazinethiol (ring- ¹³ C ₃ , 99%) | *C ₃ C ₅ H ₁₅ N ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8318-1.2 | Atrazinethiol (unlabeled) | C ₈ H ₁₅ N ₅ S | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9399-1.2 | Azinphos-methyl (unlabeled) | C ₁₀ PN ₃ H ₁₂ S ₂ O ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7140 | Bendiocarb (¹³ C ₃ , 99%) | *C ₃ C ₈ H ₁₃ NO ₄ | Inquire | |
| ULM-8638 | Bendiocarb (unlabeled) | C ₁₁ H ₁₃ NO ₄ | Inquire | |
| DLM-7152 | Bensulide (isoproxy-D ₁₄ , 98%) | C ₁₄ D ₁₄ H ₁₀ NO ₄ PS ₃ | Inquire | |
| CLM-2482-1.2 | α-HCH (α-BHC) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7232-1.2 | α-HCH (α-BHC) (unlabeled) | C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3623-1.2 | β-HCH (β-BHC) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 50 µg/mL in nonane | 2 × 1.2 mL |
| ULM-6132-1.2 | β-HCH (β-BHC) (unlabeled) | C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 2 × 1.2 mL |
| CDLM-624-1.2 | γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%; D ₆ , 99%) | *C ₆ D ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1282-1.2 | γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6133-1.2 | γ-HCH (γ-BHC) (lindane) (unlabeled) | C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3648-1.2 | δ-HCH (δ-BHC) (¹³ C ₆ , 99%) | *C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7233-1.2 | δ-HCH (δ-BHC) (unlabeled) | C ₆ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3741-1.2 | Bromoxynil (ring- ¹³ C ₆ , 99%) | *C ₆ CH ₃ Br ₂ NO | 50 µg/mL in nonane | 2 × 1.2 mL |
| ULM-6205-1.2 | Bromoxynil (unlabeled) | C ₇ H ₃ Br ₂ NO | 50 µg/mL in nonane | 1.2 mL |
| CLM-4682-1.2 | Carbaryl (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₁ NO ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-8096-1.2 | Carbaryl (unlabeled) | C ₁₀ H ₇ CO ₂ NHCH ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1911-1.2 | Carbofuran (ring- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₅ NO ₃ | 100 µg/mL in 1,4-dioxane | 1.2 mL |
| ULM-7419-1.2 | Carbofuran (unlabeled) | C ₁₂ H ₁₅ NO ₃ | 100 µg/mL in 1,4-dioxane | 1.2 mL |
| ULM-6875-1.2 | Carbofuran phenol (unlabeled) | C ₁₀ H ₁₂ O ₂ | 200 µg/mL in nonane | 1.2 mL |
| CLM-8087-1.2 | cis-Chlordane (α) (¹³ C ₁₀ , 99%) | *C ₁₀ H ₆ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-2419-1.2 | cis-Chlordane (α) (unlabeled) | C ₁₀ H ₆ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2419-25 | | | neat | 25 mg |
| CLM-4792-1.2 | trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | *C ₁₀ H ₆ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-2420-1.2 | trans-Chlordane (γ) (unlabeled) | C ₁₀ H ₆ Cl ₈ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2420-25 | | | neat | 25 mg |
| CLM-4814-1.2 | Chlordecone (kepone) (¹³ C ₁₀ , 99%) | *C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-1.2 | Chlordecone (kepone) (unlabeled) | C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-0.1 | | | neat | 0.1 g |
| CLM-4758-1.2 | Chlordene (¹³ C ₁₀ , 99%) | *C ₁₀ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7443-1.2 | Chlordene (unlabeled) | C ₁₀ H ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| CLM-6759 | 4-Chloro-2-hydroxymethyl phenoxyacetic acid (HMCPA) (ring- ¹³ C ₆ , 99%) | ClOH*C ₆ H ₃ OCH ₂ CO ₂ H | | Inquire |

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| Catalog No. | Compound | Formula | Concentration | Amount |
|--------------------------|--|---|---|---------|
| CLM-6758 | 4-Chloro-2-methylphenoxyacetic acid (MCPA) (ring- ¹³ C ₆ , 99%) | ClCH ₃ *C ₆ H ₃ OCH ₂ CO ₂ H | | Inquire |
| NEW CLM-9598-1.2 | 6-Chloronicotinic acid (¹³ C ₆ , 99%) | *C ₆ H ₄ CINO ₂ | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-9604-1.2 | 6-Chloronicotinic acid (unlabeled) | C ₆ H ₄ CINO ₂ | 100 µg/mL in MTBE | 1.2 mL |
| CLM-1913-1.2 | 4-Chlorophenol (¹³ C ₆ , 99%) | *C ₆ H ₄ CIOH | 100 µg/mL in nonane | 1.2 mL |
| ULM-7420-1.2 | 4-Chlorophenol (unlabeled) | C ₆ H ₄ CIOH | 100 µg/mL in nonane | 1.2 mL |
| DLM-4360-1.2 | Chlorpyrifos (diethyl-D ₁₀ , 99%) | C ₉ D ₁₀ HCl ₃ NO ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| ULM-7489-1.2 | Chlorpyrifos (unlabeled) | C ₉ H ₁₁ Cl ₃ NO ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-7153-1.2 | Chlorpyrifos-methyl (dimethyl-D ₆ , 98%) | C ₇ HCl ₃ D ₆ NO ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9538-1.2 | Chlorpyrifos-methyl (unlabeled) | C ₇ H ₇ Cl ₃ NO ₃ PS | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-3760-1.2 | Chlortoluron (<i>N,N</i> -dimethyl-D ₆ , 98%) | C ₁₀ H ₇ D ₆ CIN ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-3760-0.01 | | | neat | 10 mg |
| NEW ULM-9825-1.2 | Chlortoluron (unlabeled) | C ₁₀ H ₁₃ CIN ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CNLM-9940-1.2 | Clothianidin (thiazole- ¹³ C ₃ , 99%; ¹⁵ N, 98%) | *C ₃ C ₃ H ₈ Cl*NN ₄ O ₂ S | | Inquire |
| NEW ULM-9941-1.2 | Clothianidin (unlabeled) | C ₆ H ₈ CIN ₅ O ₂ S | | Inquire |
| ERC-034 | Cyclohexyl methylphosphonic acid (unlabeled) | CH ₃ PO ₃ HC ₆ H ₁₁ | 1000 µg/mL in methanol | 1.2 mL |
| CLM-7293-1.2 | Cyfluthrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7454-1.2 | Cyfluthrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₈ Cl ₂ FNO ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7292-1.2 | Cypermethrin (mix of stereoisomers) (phenoxy- ¹³ C ₆ , 99%) | *C ₆ C ₁₆ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7453-1.2 | Cypermethrin (mix of stereoisomers) (unlabeled) | C ₂₂ H ₁₉ Cl ₂ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| NEW CDLM-9205-1.2 | cis-DCCA (1,carboxyl- ¹³ C ₂ , 99%; 1-D, 97%) | C ₆ *C ₂ H ₉ DCl ₂ O ₂ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| NEW ULM-9176-1.2 | cis-DCCA (unlabeled) | C ₈ H ₁₀ Cl ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW CDLM-9206-1.2 | trans-DCCA (1,carboxyl- ¹³ C ₂ , 99%; 1-D, 97%) | C ₆ *C ₂ H ₉ DCl ₂ O ₂ | 100 µg/mL in acetonitrile-D ₃ | 1.2 mL |
| NEW ULM-9175-1.2 | trans-DCCA (unlabeled) | C ₈ H ₁₀ Cl ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-6999-1.2 | 2,4'-DDD (ring- ¹³ C ₁₂ , 99%) [(o,p'-Dichlorodiphenyl) dichloroethane] | *C ₁₂ C ₂ H ₁₀ Cl ₄ | 50 µg/mL in nonane | 1.2 mL |
| ULM-7450-1.2 | 2,4'-DDD (unlabeled) [(o,p'-Dichlorodiphenyl) dichloroethane] | C ₁₄ H ₁₀ Cl ₄ | 50 µg/mL in nonane | 1.2 mL |
| CLM-7100-1.2 | 4,4'-DDD (ring- ¹³ C ₁₂ , 99%) [(o,p'-Dichlorodiphenyl) dichloroethane] | *C ₁₂ C ₂ H ₁₀ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| DLM-3533-1.2 | 4,4'-DDD (ring-D ₈ , 98%) [(p,p'-Dichlorodiphenyl) dichloroethane] | C ₁₄ D ₈ H ₂ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7216-1.2 | 4,4'-DDD (unlabeled) [(p,p'-Dichlorodiphenyl) dichloroethane] | C ₁₄ H ₁₀ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4693-1.2 | 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) [(o,p'-Dichlorodiphenyl) dichloroethylene] | (Cl*C ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6251-1.2 | 2,4'-DDE (unlabeled) [(o,p'-Dichlorodiphenyl) dichloroethylene] | C ₁₄ H ₈ Cl ₄ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1627-1.2 | 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) [(p,p'-Dichlorodiphenyl) dichloroethylene] | (Cl*C ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1627-5 | | | neat | 5 mg |
| ULM-6137-1.2 | 4,4'-DDE (unlabeled) [(p,p'-Dichlorodiphenyl) dichloroethylene] | (ClC ₆ H ₄) ₂ C=CCl ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4692-1.2 | 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) [(o,p'-Dichlorodiphenyl) trichloroethane] | (Cl*C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6134-1.2 | 2,4'-DDT (unlabeled) [(o,p'-Dichlorodiphenyl) trichloroethane] | ClC ₆ H ₄ CH(CCl ₃)C ₆ H ₄ Cl | 100 µg/mL in nonane | 1.2 mL |
| CLM-1281-1.2 | 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) [(p,p'-Dichlorodiphenyl) trichloroethane] | (Cl*C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-1281-5 | | | neat | 5 mg |
| ULM-6135-1.2 | 4,4'-DDT (unlabeled) [(p,p'-Dichlorodiphenyl) trichloroethane] | (ClC ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-8313-1.2 | Desethylatrazine (ring- ¹³ C ₃ , 99%) CP 97% | *C ₃ C ₃ H ₁₀ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8320-1.2 | Desethylatrazine (unlabeled) | C ₆ H ₁₀ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-7528-1.2 | Desethyl desisopropyl atrazine (¹³ C ₃ , 99%) CP 95% | *C ₃ H ₄ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8001-1.2 | Desethyl desisopropyl atrazine (unlabeled) | C ₃ H ₄ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-8315-1.2 | Desethylhydroxyatrazine (ring- ¹³ C ₃ , 99%) | C ₃ *C ₃ H ₁₁ N ₅ O | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |

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| Catalog No. | Compound | Formula | Concentration | Amount | |
|--------------------|--|--|--|---------------------------|--------|
| ULM-8322-1.2 | Desethylhydroxyatrazine (unlabeled) | C ₆ H ₁₁ N ₅ O | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL | |
| CLM-8312-1.2 | Desisopropylatrazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₂ H ₈ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL | |
| ULM-8319-1.2 | Desisopropylatrazine (unlabeled) | C ₅ H ₈ CIN ₅ | 100 µg/mL in acetonitrile | 1.2 mL | |
| CLM-8314-1.2 | Desisopropylhydroxyatrazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₂ H ₉ N ₅ O | 100 µg/mL in acetonitrile | 1.2 mL | |
| ULM-8321-1.2 | Desisopropylhydroxyatrazine (unlabeled) | C ₅ H ₉ N ₅ O | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL | |
| DLM-1148-1.2 | Diazinon (diethyl-D ₁₀ , 98%) | C ₁₂ H ₁₁ D ₁₀ N ₂ O ₃ PS | 100 µg/mL in nonane | 1.2 mL | |
| NEW | DLM-1148-A-1.2 | | 100 µg/mL in acetonitrile | 1.2 mL | |
| | DLM-1148-5 | | neat | 5 mg | |
| ULM-6575-S-10X-1.2 | Diazinon (unlabeled) | C ₁₂ H ₂₁ N ₂ O ₃ PS | 1000 µg/mL in nonane | 1.2 mL | |
| NEW | ULM-6575-A-1.2 | | 100 µg/mL in acetonitrile | 1.2 mL | |
| CLM-816-1.2 | 2,6-Dichloro-4-nitroaniline (ring- ¹³ C ₆ , 99%) | Cl ₂ *C ₆ H ₄ (NO ₂)NH ₂ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-1858-1.2 | 2,4-Dichlorophenoxyacetic acid (2,4-D) (ring- ¹³ C ₆ , 99%) | Cl ₂ *C ₆ H ₃ OCH ₂ CO ₂ H | 100 µg/mL in acetonitrile | 1.2 mL | |
| DLM-1146-5 | 2,4-Dichlorophenoxyacetic acid (2,4-D) (ring-D ₃ , 98%) | Cl ₂ C ₆ D ₃ OCH ₂ CO ₂ H | neat | 5 mg | |
| ULM-7418-1.2 | 2,4-Dichlorophenoxyacetic acid (2,4-D) (unlabeled) | Cl ₂ C ₆ H ₃ OCH ₂ CO ₂ H | 100 µg/mL in acetonitrile | 1.2 mL | |
| CLM-3722-1.2 | Dichlorprop (ring- ¹³ C ₆ , 99%) | *C ₆ C ₃ H ₈ Cl ₂ O ₃ | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7313-1.2 | Dichlorprop (unlabeled) | (Cl) ₂ C ₆ H ₃ OCH(CH ₃)CO ₂ H | 100 µg/mL in nonane | 1.2 mL | |
| DLM-2829-0.01 | Dichlorvos (dimethyl-D ₆ , 98%) | C ₄ D ₆ Cl ₂ O ₄ P | neat | 10 mg | |
| ULM-7217-1.2 | Dichlorvos (unlabeled) | (H ₃ CO) ₂ POOCH=CCl ₂ | 100 µg/mL in nonane | 1.2 mL | |
| CLM-4726-1.2 | Dieldrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL | |
| ULM-7230-1.2 | Dieldrin (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL | |
| NEW | DLM-4851-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| | ULM-9287-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) (unlabeled) | C ₄ H ₁₀ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| DLM-4852-1.2 | O,O-Diethyl thiophosphate, potassium salt (DETP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₃ PS | 100 µg/mL in methanol | 1.2 mL | |
| ERD-119 | O,O-Diethyl thiophosphate, potassium salt (DETP) (unlabeled) | C ₄ H ₁₀ KO ₃ PS | 1000 µg/mL in methanol | 1.2 mL | |
| NEW | DLM-9003-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| | ULM-9002-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) (unlabeled) | C ₄ H ₁₀ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| DLM-4762-1.2 | N,N-Diethyl-m-toluamide (DEET) (dimethyl-D ₆ , 98%) | CH ₃ C ₆ H ₄ CON(CH ₂ CD ₃) ₂ | 100 µg/mL in MeCl-D ₂ | 1.2 mL | |
| DLM-4762-D-1.2 | | | 100 µg/mL in dioxane | 1.2 mL | |
| ULM-7975-1.2 | N,N-Diethyl-m-toluamide (DEET) (unlabeled) | CH ₃ C ₆ H ₄ CON(CH ₂ CH ₃) ₂ | 100 µg/mL in MeCl | 1.2 mL | |
| ULM-7975-D-1.2 | | | 100 µg/mL in dioxane | 1.2 mL | |
| NEW | ULM-9898-1.2 | Diisopropyl methylphosphonate (unlabeled) | C ₇ H ₁₇ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| | DLM-7151-1.2 | Dimethoate (O,O-dimethyl-D ₆ , 98%) | C ₅ D ₆ H ₆ NO ₃ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7972-1.2 | Dimethoate (unlabeled) | C ₅ H ₁₂ NO ₃ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL | |
| NEW | DLM-8868-1.2 | O,O-Dimethyl phosphate, potassium salt (DMP) (dimethyl-D ₆ , 98%) | C ₂ D ₆ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| | ULM-8867-1.2 | O,O-Dimethyl phosphate, potassium salt (DMP) (unlabeled) | C ₂ H ₆ KO ₃ P ₄ | 100 µg/mL in methanol | 1.2 mL |
| NEW | DLM-8904-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) (dimethyl-D ₆ , 98%) CP 97% | C ₂ D ₆ KO ₃ PS | 100 µg/mL in methanol | 1.2 mL |
| | ULM-8905-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) (unlabeled) CP 97% | C ₂ H ₆ KO ₃ PS | 1000 µg/mL in methanol | 1.2 mL |
| NEW | DLM-4541-M-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) (dimethyl-D ₆ , 98%) | C ₂ D ₆ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| | ULM-9004-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) (unlabeled) | C ₂ H ₆ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| CLM-3373 | Dinocap (ring- ¹³ C ₆ , 99%) | *C ₆ C ₁₂ H ₂₄ N ₂ O ₆ | Inquire | | |
| NEW | ULM-9171-1.2 | Dinocap (unlabeled) | C ₁₈ H ₂₄ N ₂ O ₆ | Inquire | |
| | CLM-9594-1.2 | Dinotefuran (furylmethyl- ¹³ C ₅ , 99%) | *C ₅ C ₂ H ₁₄ N ₄ O ₃ | 100 µg/mL in methanol | 1.2 mL |
| ULM-9732-1.2 | Dinotefuran (unlabeled) | C ₇ H ₁₄ N ₄ O ₃ | 100 µg/mL in methanol | 1.2 mL | |

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| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------------------|--|---|---------------------------|---------|
| NEW ULM-9899-1.2 | Dipinacolyl methylphosphonate (unlabeled) | C ₁₃ H ₂₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| DLM-7183 | Disulfoton (<i>O,O</i> -diethyl-D ₁₀ , 98%) | C ₈ D ₁₀ H ₉ O ₂ PS ₃ | | Inquire |
| CLM-6025-1.2 | Endosulfan I (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| DLM-2862-1.2 | Endosulfan I (D ₄ , 97%) | C ₉ D ₄ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7447-1.2 | Endosulfan I (unlabeled) | C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-6026-1.2 | Endosulfan II (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7448-1.2 | Endosulfan II (unlabeled) | C ₉ H ₆ Cl ₆ O ₃ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-7531-1.2 | Endosulfan sulfate (¹³ C ₉ , 99%) | *C ₉ H ₆ Cl ₆ O ₄ S | 100 µg/mL in nonane | 1.2 mL |
| ULM-7990-1.2 | Endosulfan sulfate (unlabeled) | C ₉ H ₆ Cl ₆ O ₄ S | 100 µg/mL in nonane | 1.2 mL |
| CLM-4782-1.2 | Endrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-7444-1.2 | Endrin (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-4815-1.2 | Endrin aldehyde (¹³ C ₁₂ , 99%) | *C ₁₂ H ₁₀ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-4815-50 | | | neat | 50 µg |
| NEW ULM-8958-1.2 | Endrin aldehyde (unlabeled) | C ₁₂ H ₁₀ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8958-50 | | | neat | 50 µg |
| NEW CLM-4816-1.2 | Endrin ketone (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-4816-50 | | | neat | 50 µg |
| NEW ULM-8956-1.2 | Endrin ketone (unlabeled) | C ₁₂ H ₈ Cl ₆ O | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-8956-50 | | | neat | 50 µg |
| NEW ULM-6091-1.2 | Ethyl dimethylamidophosphate, sodium salt (unlabeled) | C ₄ H ₁₁ NO ₃ PNa | 1000 µg/mL in methanol | 1.2 mL |
| NEW DLM-6098-1.2 | Ethyl methylphosphonate (ethyl-D ₅ , 98%) | C ₃ H ₄ D ₅ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-6099-1.2 | Ethyl methylphosphonic acid (unlabeled) | C ₃ H ₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| DLM-2878-0.01 | Fenitrothion (<i>O,O</i> -dimethyl-D ₆ , 98%) | C ₉ D ₆ H ₆ NO ₅ PS | neat | 10 mg |
| NEW CNLM-9636-1.2 | Fipronil (3-cyano, pyrazole-3,4,5- ¹³ C ₄ , 99%; 3-cyano, 5- ¹⁵ N ₂ , 98%) | *C ₄ C ₈ H ₄ Cl ₂ F ₆ *N ₂ OS | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9635-1.2 | Fipronil (unlabeled) | C ₁₂ H ₄ Cl ₂ F ₆ N ₄ OS | 100 µg/mL in methanol | 1.2 mL |
| NEW CNLM-9647-1.2 | Fipronil desulfanyl (3-cyano, pyrazole-3,4,5- ¹³ C ₄ , 99%; *C ₄ C ₈ H ₄ Cl ₂ F ₆ *N ₂) 3-cyano, 5- ¹⁵ N ₂ , 98%) | | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9646-1.2 | Fipronil desulfanyl (unlabeled) | C ₁₂ H ₄ Cl ₂ F ₆ N ₄ | 100 µg/mL in methanol | 1.2 mL |
| NEW CNLM-9650-1.2 | Fipronil difluoromethyl sulfinyl (3-cyano, pyrazole-3,4,5- ¹³ C ₄ , 99%) 3-cyano- 5- ¹⁵ N ₂ , 98%) | *C ₄ C ₇ H ₅ Cl ₂ F ₃ *N ₂ N ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9649-1.2 | Fipronil trifluoromethyl sulfinyl (unlabeled) | C ₁₁ H ₅ Cl ₂ F ₃ N ₂ N ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW CNLM-9645-1.2 | Fipronil sulfide (3-cyano, pyrazole-3,4,5- ¹³ C ₄ , 99%; 3-cyano, 5- ¹⁵ N ₂ , 98%) | *C ₄ C ₈ H ₄ Cl ₂ F ₆ *N ₂ N ₂ S | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9644-1.2 | Fipronil sulfide (unlabeled) | C ₁₂ H ₄ Cl ₂ F ₆ N ₄ S | 100 µg/mL in methanol | 1.2 mL |
| NEW CNLM-9643-1.2 | Fipronil sulfone (3-cyano, pyrazole-3,4,5- ¹³ C ₄ , 99%; *C ₄ C ₈ H ₄ Cl ₂ F ₆ *N ₂ N ₂ O ₂ S | | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9642-1.2 | Fipronil sulfone (unlabeled) | C ₁₂ H ₄ Cl ₂ F ₆ N ₄ O ₂ S | 100 µg/mL in methanol | 1.2 mL |
| CLM-7389-1.2 | 4-Fluoro-3-phenoxybenzoic acid (¹³ C ₆ , 99%) | *C ₆ C ₇ H ₆ FO ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7391-1.2 | 4-Fluoro-3-phenoxybenzoic acid (unlabeled) | C ₁₃ H ₉ FO ₃ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-4545-1.2 | Fonofos (ring- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₁₅ OPS ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6694-1.2 | Fonofos (unlabeled) | CH ₂ CH ₃ P(S)(OCH ₂ CH ₃)(SC ₆ H ₅) | 100 µg/mL in nonane | 1.2 mL |
| NEW DLM-9680 | Forchlorfuron (phenyl-D ₅ , 99%) | C ₁₂ D ₅ H ₅ ClN ₃ O | | Inquire |
| CNLM-4666-1.2 | Glyphosate (2- ¹³ C, 99%; ¹⁵ N, 98%+) | HO ₂ *CCH ₂ *NHCH ₂ PO(OH) ₂ | 100 µg/mL in water | 1.2 mL |
| NEW CNLM-4666-10X-1.2 | | | 1000 µg/mL in water | 1.2 mL |
| CNLM-4666-10 | | | 100 µg/mL in water | 10 mL |
| ULM-6876-1.2 | Glyphosate (unlabeled) | HO ₂ CCH ₂ NHCH ₂ PO(OH) ₂ | 100 µg/mL in water | 1.2 mL |
| CLM-4759-1.2 | Heptachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₇ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-1.2 | Heptachlor (unlabeled) | C ₁₀ H ₅ Cl ₇ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2424-0.1 | | | neat | 0.1 g |
| CLM-4734-1.2 | cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₇ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2425-1.2 | cis-Heptachlor epoxide (unlabeled) | C ₁₀ H ₅ Cl ₇ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2425-0.1 | | | neat | 0.1 g |
| ULM-7869-1.2 | trans-Heptachlor epoxide (unlabeled) | C ₁₀ H ₅ Cl ₇ O | 100 µg/mL in nonane | 1.2 mL |
| CLM-351-1.2 | Hexachlorobenzene (¹³ C ₆ , 99%) | *C ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-6130-1.2 | Hexachlorobenzene (unlabeled) | C ₆ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |

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Individual Pesticide and Pesticide Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|---|---|---|--------|
| NEW ULM-9429-1.2 | Hp-Sed (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9428-1.2 | Hx-Sed (unlabeled) | C ₁₀ H ₁₂ Cl ₆ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8310-1.2 | Hydroxyatrazine (ring- ¹³ C ₃ , 99%) | (CH ₃ CH ₂ NH)*C ₃ N ₃ (OH) (NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| ULM-8317-1.2 | Hydroxyatrazine (unlabeled) | (CH ₃ CH ₂ NH)C ₃ N ₃ (OH) (NHCH(CH ₃) ₂) | 100 µg/mL in 80% water/ 20% diethylamine | 1.2 mL |
| DLM-8512-1.2 | Imidacloprid (4,4,5,5-D ₄ , 98%) | C ₉ H ₆ D ₄ ClN ₅ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| ULM-8513-1.2 | Imidacloprid (unlabeled) | C ₉ H ₁₀ ClN ₅ O ₂ | 100 µg/mL in methanol | 1.2 mL |
| CLM-4727-1.2 | Isodrin (¹³ C ₁₂ , 99%) | *C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7442-1.2 | Isodrin (unlabeled) | C ₁₂ H ₈ Cl ₆ | 100 µg/mL in nonane | 1.2 mL |
| ERI-026 | Isobutyl hydrogen methylphosphonate (unlabeled) | C ₅ H ₁₃ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| ERI-015 | Isopropyl methylphosphonic acid (unlabeled) | C ₄ H ₁₁ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| CLM-4814-1.2 | Kepone (chlordecone) (¹³ C ₁₀ , 99%) | *C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-1.2 | Kepone (chlordecone) (unlabeled) | C ₁₀ Cl ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-2301-0.1 | | | neat | 0.1 g |
| DLM-4476-1.2 | Malathion (D ₁₀ , 99%) | C ₁₀ D ₁₀ H ₉ O ₆ PS ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-8122-1.2 | Malathion (unlabeled) | C ₁₀ H ₁₉ O ₆ PS ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW CLM-9050-1.2 | Malathion diacid (¹³ C ₄ , 99%) CP 97% | *C ₄ C ₂ H ₁₁ O ₆ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9073-1.2 | Malathion diacid (unlabeled) | C ₆ H ₁₁ O ₆ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW DLM-7149-1.2 | Methamidophos (dimethyl-D ₆ , 98%) | C ₂ D ₆ H ₂ NO ₂ PS | 100 µg/mL in dioxane | 1.2 mL |
| NEW ULM-8872-1.2 | Methamidophos (unlabeled) | C ₂ H ₈ NO ₂ PS | 100 µg/mL in dioxane | 1.2 mL |
| CNLM-7148-1.2 | Methomyl (acetohydroxamate- ¹³ C ₂ , 99%; ¹⁵ N, 98%) | *C ₂ C ₃ H ₁₀ N*NO ₂ S | 100 µg/mL in methanol | 1.2 mL |
| ULM-8639-1.2 | Methomyl (unlabeled) | C ₅ H ₁₀ NNO ₂ S | 100 µg/mL in methanol | 1.2 mL |
| CLM-4683-1.2 | Methoxychlor (ring- ¹³ C ₁₂ , 99%) | (CH ₃ O)*C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7440-1.2 | Methoxychlor (unlabeled) | (CH ₃ O)C ₆ H ₄) ₂ CHCCl ₃ | 100 µg/mL in nonane | 1.2 mL |
| CDLM-6100-1.2 | Methylphosphonic acid (¹³ C, 99%; methyl-D ₃ , 98%) | *CD ₃ H ₂ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| DLM-6196-1.2 | Methylphosphonic acid (methyl-D ₃ , 98%) | CD ₃ P(O)(OH) ₂ | 100 µg/mL in methanol | 1.2 mL |
| ERM-038 | Methylphosphonic acid (unlabeled) | CH ₃ P(O)(OH) ₂ | 100 µg/mL in methanol | 1.2 mL |
| CLM-6620-1.2 | Methylphosphonic acid, mono-(1,2,2-trimethylpropyl) ester (trimethylpropyl- ¹³ C ₆ , 99%) | *C ₆ CH ₁₇ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| CLM-3712-1.2 | Metolachlor (ring- ¹³ C ₆ , 99%) | *C ₆ C ₉ H ₂₂ ClNO ₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7314-1.2 | Metolachlor (unlabeled) | C ₁₅ H ₂₂ ClNO ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4813-1.2 | Mirex (¹³ C ₁₀ , 99%) | *C ₁₀ Cl ₁ | 100 µg/mL in nonane | 1.2 mL |
| CLM-2078-1 | Mirex (¹³ C ₈ , 99%) | *C ₈ C ₂ Cl ₁₂ | 200 µg/mL in toluene | 1 mL |
| ULM-2427-1.2 | Mirex (unlabeled) | C ₁₀ Cl ₁₂ | 100 µg/mL in nonane | 1.2 mL |
| ULM-2427-0.1 | | | neat | 0.1 g |
| CLM-4811-1.2 | cis-Nonachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7445-1.2 | cis-Nonachlor (unlabeled) | C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4735-1.2 | trans-Nonachlor (¹³ C ₁₀ , 99%) | *C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7229-1.2 | trans-Nonachlor (unlabeled) | C ₁₀ H ₅ Cl ₉ | 100 µg/mL in nonane | 1.2 mL |
| CLM-4729-1.2 | Oxychlordane (¹³ C ₁₀ , 99%) | *C ₁₀ H ₄ Cl ₈ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-1.2 | Oxychlordane (unlabeled) | C ₁₀ H ₄ Cl ₈ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-6139-SM-1.2 | | | 100 µg/mL in methanol | 1.2 mL |
| DLM-7150-1.2 | Oxydemeton methyl (O,O-dimethyl-D ₆ , 98%) | C ₆ D ₆ H ₉ O ₄ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8579-1.2 | Oxydemeton methyl (unlabeled) | C ₆ H ₁₅ O ₄ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-4538-1.2 | Oxypyrimidine (diazinon metabolite) (methyl-4,5,6- ¹³ C ₄ , 99%) | *C ₄ C ₄ H ₁₂ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7432-1.2 | Oxypyrimidine (unlabeled) | C ₈ H ₁₂ N ₂ O | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-2970-1.2 | Parathion (diethyl-D ₁₀ , 98%) | C ₁₀ D ₁₀ H ₄ NO ₅ PS | 100 µg/mL in nonane | 1.2 mL |
| ULM-8144-1.2 | Parathion (unlabeled) | NO ₂ (C ₆ H ₄)OP(=S)(OC ₂ H ₅) ₂ | 100 µg/mL in nonane | 1.2 mL |
| CLM-7930-1.2 | Parlar 26 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7828-1.2 | Parlar 26 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8705-1.2 | Parlar 32 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8665-1.2 | Parlar 32 (unlabeled) | C ₁₀ H ₁₁ Cl ₇ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9005-1.2 | Parlar 38 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8719-1.2 | Parlar 39 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8767-1.2 | Parlar 39 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9430-1.2 | Parlar 40 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9431-1.2 | Parlar 41 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |
| NEW ULM-9432-1.2 | Parlar 44 (unlabeled) | C ₁₀ H ₁₀ Cl ₈ | 10 µg/mL in nonane | 1.2 mL |

Individual Pesticide and Pesticide Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|---------------------------|--|--|---------------------------|--------|
| CLM-7931-1.2 | Parlar 50 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7829-1.2 | Parlar 50 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7932-1.2 | Parlar 62 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-7830-1.2 | Parlar 62 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8720-1.2 | Parlar 69 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8768-1.2 | Parlar 69 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-8721-1.2 | Parlar 70 (¹³ C ₁₀ , 99%) | *C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| ULM-8769-1.2 | Parlar 70 (unlabeled) | C ₁₀ H ₉ Cl ₉ | 10 µg/mL in nonane | 1.2 mL |
| CLM-7322-1.2 | cis-Permethrin (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8526-1.2 | cis-Permethrin (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| CLM-7323-1.2 | trans-Permethrin (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| ULM-8527-1.2 | trans-Permethrin (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CH ₂ CO ₂ C ₇ H ₉ Cl ₂ O ₃ | 50 µg/mL in nonane | 1.2 mL |
| CLM-4542-1.2 | 3-Phenoxybenzoic acid (phenoxy- ¹³ C ₆ , 99%) | *C ₆ H ₅ OC ₆ H ₄ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |
| CLM-4542-SA-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-6781-1.2 | 3-Phenoxybenzoic acid (unlabeled) | C ₆ H ₅ OC ₆ H ₄ CO ₂ H | 100 µg/mL in nonane | 1.2 mL |
| ULM-6781-SA-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-3733-1.2 | o-Phenylphenol (phenyl- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₀ O | 100 µg/mL in nonane | 1.2 mL |
| ULM-7396-1.2 | o-Phenylphenol (unlabeled) | C ₁₂ H ₉ OH | 100 µg/mL in nonane | 1.2 mL |
| CLM-3748-1.2 | p-Phenylphenol (phenyl- ¹³ C ₆ , 99%) | *C ₆ C ₆ H ₁₀ O | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-4544-1.2 | Phorate (diethoxy- ¹³ C ₄ , 99%) | (*C ₂ H ₅ O) ₂ P(S)SCH ₂ SC ₂ H ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7567-1.2 | Phorate (unlabeled) | (C ₂ H ₅ O) ₂ P(S)SCH ₂ SC ₂ H ₅ | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-4667-1.2 | Phosmet (dimethyl-D ₆ , 98%) | C ₁₁ H ₆ D ₆ NO ₄ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-8454-1.2 | Phosmet (unlabeled) | C ₁₁ H ₁₂ NO ₄ PS ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| CLM-3738-1.2 | Propazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₆ H ₁₆ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8304 | Propazine (unlabeled) | C ₉ H ₁₆ CIN ₅ | Inquire | |
| ULM-7141-1.2 | Propoxur (isopropyl-D ₇ , 98%) | C ₁₁ D ₇ H ₈ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| NEW ULM-9765-1.2 | Propoxur (unlabeled) | C ₁₁ H ₁₅ NO ₃ | 100 µg/mL in nonane | 1.2 mL |
| CLM-3739-1.2 | Simazine (ring- ¹³ C ₃ , 99%) | *C ₃ C ₄ H ₁₂ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW CLM-3739-A-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| ULM-7893-1.2 | Simazine (unlabeled) | C ₇ H ₁₂ CIN ₅ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-7893-A-1.2 | | | 100 µg/mL in acetonitrile | 1.2 mL |
| CDLM-7943-1.2 | Sodium monofluoroacetate (¹³ C ₂ , 99%; 2,2-D ₂ , 98%) | F*CD ₂ *CO ₂ Na | 1 mg/mL in water | 1.2 mL |
| NEW ULM-9674-1.2 | Sodium monofluoroacetate (unlabeled) | FCH ₂ CO ₂ Na | 1 mg/mL in water | 1.2 mL |
| DLM-380-1.2 | Styrene (D ₈ , 98%) + BHT | C ₆ D ₅ CD=CD ₂ | 100 µg/mL in nonane | 1.2 mL |
| NEW CNLM-9869-1.2 | Sulfoxaflor (cyano- ¹³ C, 99%; cyano- ¹⁵ N, imine- ¹⁵ N, 98%) | *CC ₉ H ₁₀ F ₃ *N ₂ NOS | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9870-1.2 | Sulfoxaflor (unlabeled) | C ₁₀ H ₁₀ F ₃ N ₂ NOS | 100 µg/mL in methanol | 1.2 mL |
| CLM-4543 | Terbufos (diethoxy- ¹³ C ₄ , 99%) | C(CH ₃) ₃ SCH ₂ SP(S)(O*CH ₂ *CH ₃) ₂ | Inquire | |
| NEW CLM-9690-1.2 | 3-Tetrahydrofuroic acid (¹³ C ₅ , 99%) | *C ₅ H ₈ O ₃ | 100 µg/mL in MTBE | 1.2 mL |
| NEW ULM-9691-1.2 | 3-Tetrahydrofuroic acid (unlabeled) | C ₅ H ₈ O ₃ | 100 µg/mL in MTBE | 1.2 mL |
| NEW CLM-9652-1.2 | Thiacloprid (pyridylmethyl- ¹³ C ₆ , 99%) | *C ₆ C ₄ H ₉ CIN ₄ S | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9733-1.2 | Thiacloprid (unlabeled) | C ₁₀ H ₉ CIN ₄ S | 100 µg/mL in methanol | 1.2 mL |
| NEW CNLM-9860-1.2 | Thiamethoxam (thiazole- ¹³ C ₃ , 99%; ¹⁵ N, 98%) | *C ₃ C ₅ H ₁₀ Cl*NN ₄ O ₃ S | Inquire | |
| NEW ULM-9939-1.2 | Thiamethoxam (unlabeled) | C ₈ H ₁₆ CIN ₃ O ₃ S | Inquire | |
| CLM-4551-1.2 | 2,4,5-Trichlorophenoxyacetic acid (2,4,5-T) (ring- ¹³ C ₆ , 99%) | *C ₆ C ₂ H ₅ Cl ₃ O ₃ | 100 µg/mL in MeCl | 1.2 mL |
| ULM-7213-1.2 | 2,4,5-Trichlorophenoxyacetic acid (2,4,5-T) (unlabeled) | C ₆ H ₂ Cl ₃ OCH ₂ CO ₂ H | 100 µg/mL in MeCl | 1.2 mL |
| NEW CLM-9049-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (4,5,6- ¹³ C ₃ , 99%) CP 97% | *C ₃ C ₂ H ₂ Cl ₃ NO | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9204-1.2 | 3,5,6-Trichloro-2-pyridinol (TCPY) (unlabeled) | C ₅ H ₂ Cl ₃ NO | 100 µg/mL in acetonitrile | 1.2 mL |
| DLM-4479-1.2 | Trifluralin (di-n-propyl-D ₁₄ , 98%) | C ₁₃ D ₁₄ H ₂ F ₃ N ₃ O ₄ | 100 µg/mL in nonane | 1.2 mL |
| ULM-7236-1.2 | Trifluralin (unlabeled) | C ₁₃ H ₁₄ H ₂ F ₃ N ₃ O ₄ | 100 µg/mL in nonane | 1.2 mL |

Stockholm Convention POPs Pesticide Standard Mixtures

| Catalog No. | Compound | Amount | | | | | |
|---|---|--|-----|-----|-----|-----|-----|
| NEW | ES-5464-A | Expanded POPs Pesticides Calibration Solutions with Endosulfan Sulfate [CS1-CS6] | | | | | |
| Individual calibration solutions are available. Please inquire. | | | | | | | |
| | | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
| | Unlabeled | | | | | | |
| | Hexachlorobenzene | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Pentachlorobenzene | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Aldrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Dieldrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Endrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 4,4'-DDT | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 4,4'-DDE | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 4,4'-DDD | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 2,4'-DDT | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 2,4'-DDE | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | 2,4'-DDD | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>trans</i> -Chlordane (γ) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>cis</i> -Chlordane (α) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>trans</i> -Nonachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>cis</i> -Nonachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Oxychlordane | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Heptachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>trans</i> -Heptachlor epoxide | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | <i>cis</i> -Heptachlor epoxide | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Mirex | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Kepone (chlordecone) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | α -HCH (α -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | β -HCH (β -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | γ -HCH (γ -BHC) (lindane) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | δ -HCH (δ -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Endosulfan I | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Endosulfan II | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Endosulfan sulfate | 0.4 | 2 | 10 | 40 | 200 | 800 |
| | Labeled | | | | | | |
| | Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Pentachlorobenzene ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Aldrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Endrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Dieldrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 2,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | <i>trans</i> -Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | <i>cis</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Oxychlordane ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Heptachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Mirex ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Kepone (chlordecone) ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | δ -HCH (δ -BHC) ($^{13}\text{C}_6$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Endosulfan I ($^{13}\text{C}_9$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Endosulfan II ($^{13}\text{C}_9$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Endosulfan sulfate ($^{13}\text{C}_9$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Syringe | | | | | | |
| | 4,4'-DiCB ($^{13}\text{C}_{12}$, 99%) (PCB-15) | 20 | 20 | 20 | 20 | 20 | 20 |
| | 2,3',4',5-TetraCB ($^{13}\text{C}_{12}$, 99%) (PCB-70) | 20 | 20 | 20 | 20 | 20 | 20 |
| | Sampling | | | | | | |
| | Isodrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |

Stockholm Convention POPs Pesticide Standard Mixtures

| Catalog No. | Compound | Amount | | | | | |
|--|--|--|-----|-----|-----|-----|-----|
| ES-5464 | Expanded POPs Pesticides Calibration Solutions [CS1-CS6] | Set of 6 × 0.2 mL in nonane | | | | | |
| <i>Individual calibration solutions are available. Please inquire.</i> | | | | | | | |
| | | <i>All concentrations are in ng/mL (ppb)</i> | | | | | |
| Unlabeled | | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 |
| Hexachlorobenzene | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Pentachlorobenzene | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Aldrin | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Dieldrin | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Endrin | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDT | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDE | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDD | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDT | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDE | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDD | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Chlordane (γ) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Chlordane (α) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Nonachlor | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Nonachlor | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Oxychlordane | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Heptachlor | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Heptachlor epoxide | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Heptachlor epoxide | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Mirex | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Kepone (chlordecone) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| α -HCH (α -BHC) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| β -HCH (β -BHC) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| γ -HCH (γ -BHC) (lindane) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| δ -HCH (δ -BHC) | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Endosulfan I | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Endosulfan II | | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Labeled | | | | | | | |
| Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Pentachlorobenzene ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Aldrin ($^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Endrin ($^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Dieldrin ($^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>trans</i> -Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>cis</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Oxychlordane ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Heptachlor ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Mirex ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Kepone (chlordecone) ($^{13}\text{C}_{10}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| δ -HCH (δ -BHC) ($^{13}\text{C}_6$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Endosulfan I ($^{13}\text{C}_9$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Endosulfan II ($^{13}\text{C}_9$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Syringe | | | | | | | |
| 4,4'-DiCB ($^{13}\text{C}_{12}$, 99%) (PCB-15) | | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3',4',5-TetraCB ($^{13}\text{C}_{12}$, 99%) (PCB-70) | | 20 | 20 | 20 | 20 | 20 | 20 |
| Sampling | | | | | | | |
| Isodrin ($^{13}\text{C}_{12}$, 99%) | | 20 | 20 | 20 | 20 | 20 | 20 |

Stockholm Convention POPs Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------------------|---|------------------|
| NEW ES-5465-A | Expanded POPs Pesticides Cleanup Spike with Endosulfan Sulfate | 1.2 mL in nonane |
| NEW ES-5465-A-5X | Expanded POPs Pesticides Cleanup Spike with Endosulfan Sulfate (5X stock) | 1.2 mL in nonane |
| ES-5465 | Expanded POPs Pesticides Cleanup Spike | 1.2 mL in nonane |
| ES-5465-5X | Expanded POPs Pesticides Cleanup Spike (5X stock) | 1.2 mL in nonane |

| Labeled | ES-5465-A (ng/mL) | ES-5465-A-5X (ng/mL) | ES-5465 (ng/mL) | ES-5465-5X (ng/mL) |
|--|----------------------|-------------------------|--------------------|-----------------------|
| Hexachlorobenzene (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| Pentachlorobenzene (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| Aldrin (¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| Endrin (¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| Dieldrin (¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 4,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| 2,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 | 500 | 100 | 500 |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| cis-Nonachlor (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| Heptachlor (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| Mirex (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| Kepone (chlordecone) (¹³ C ₁₀ , 99%) | 100 | 500 | 100 | 500 |
| α -HCH (α -BHC) (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| β -HCH (β -BHC) (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| γ -HCH (γ -BHC) (lindane) (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| δ -HCH (δ -BHC) (¹³ C ₆ , 99%) | 100 | 500 | 100 | 500 |
| Endosulfan I (¹³ C ₉ , 99%) | 100 | 500 | 100 | 500 |
| Endosulfan II (¹³ C ₉ , 99%) | 100 | 500 | 100 | 500 |
| Endosulfan sulfate (¹³ C ₉ , 99%) | 100 | 500 | — | — |

| | | |
|-----------|--|------------------|
| EC-5350 | POPs Pesticides HRMS (PCB) Syringe Spike | 1.2 mL in nonane |
| EC-5350-L | POPs Pesticides LRMS (PCB) Syringe Spike | 0.5 mL in nonane |

| Labeled | ES-5350 (ng/mL) | ES-5350-L (ng/mL) |
|---|--------------------|----------------------|
| 4,4'-DiCB (¹³ C ₁₂ , 99%) (PCB-15) | 100 | 1000 |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) (PCB-70) | 100 | 1000 |

| | | |
|---------|---|------------------|
| ES-5466 | Expanded POPs Pesticides Sampling Spike | 1.2 mL in nonane |
|---------|---|------------------|

| Labeled | (ng/mL) |
|--|---------|
| Isodrin (¹³ C ₁₂ , 99%) | 1000 |

Stockholm Convention POPs Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|--|---|--------------------|
| NEW | | |
| ES-5467-A | Expanded POPs Pesticides PAR Solution with Endosulfan Sulfate | 1.2 mL in nonane |
| ES-5467 | Expanded POPs Pesticides PAR Solution | 1.2 mL in nonane |
| | | |
| Unlabeled | | |
| Hexachlorobenzene | ES-5467-A (ng/mL) | ES-5467 (ng/mL) |
| Pentachlorobenzene | 1000 | 1000 |
| Aldrin | 1000 | 1000 |
| Dieldrin | 1000 | 1000 |
| Endrin | 1000 | 1000 |
| 4,4'-DDT | 1000 | 1000 |
| 4,4'-DDE | 1000 | 1000 |
| 4,4'-DDD | 1000 | 1000 |
| 2,4'-DDT | 1000 | 1000 |
| 2,4'-DDE | 1000 | 1000 |
| 2,4'-DDD | 1000 | 1000 |
| <i>trans</i> -Chlordane (γ) | 1000 | 1000 |
| <i>cis</i> -Chlordane (α) | 1000 | 1000 |
| <i>trans</i> -Nonachlor | 1000 | 1000 |
| <i>cis</i> -Nonachlor | 1000 | 1000 |
| Oxychlordane | 1000 | 1000 |
| Heptachlor | 1000 | 1000 |
| <i>trans</i> -Heptachlor epoxide | 1000 | 1000 |
| <i>cis</i> -Heptachlor epoxide | 1000 | 1000 |
| Mirex | 1000 | 1000 |
| Kepone (chlordecone) | 1000 | 1000 |
| α -HCH (α -BHC) | 1000 | 1000 |
| β -HCH (β -BHC) | 1000 | 1000 |
| γ -HCH (γ -BHC) (lindane) | 1000 | 1000 |
| δ -HCH (δ -BHC) | 1000 | 1000 |
| Endosulfan I | 1000 | 1000 |
| Endosulfan II | 1000 | 1000 |
| Endosulfan sulfate | 1000 | — |

Toxaphene Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| NEW ES-5543 | US EPA Method 8276 Toxaphene Composite Stock Standard | 1.2 mL in nonane |
| Unlabeled (ng/mL) | | |
| Hx-Sed | 1000 | |
| Hp-Sed | 1000 | |
| Parlar 26 | 1000 | |
| Parlar 40 | 1000 | |
| Parlar 41 | 1000 | |
| Parlar 44 | 1000 | |
| Parlar 50 | 1000 | |
| Parlar 62 | 1000 | |
| Labeled (ng/mL) | | |
| Parlar 26 (¹³ C ₁₀ , 99%) | 100 | |
| Parlar 50 (¹³ C ₁₀ , 99%) | 100 | |
| Parlar 62 (¹³ C ₁₀ , 99%) | 100 | |
| NEW ES-5544 | US EPA Method 8276 Toxaphene Surrogate Standard | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| Parlar 39 (¹³ C ₁₀ , 99%) | 100 | |
| NEW ES-5545 | US EPA Method 8276 Toxaphene Injection Internal Standard | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 100 | |
| ES-5352-L | POPs Toxaphene Surrogate Solution with PCB Syringe | 1.2 mL in nonane |
| Labeled (ng/mL) | | |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 1000 | |
| ES-5353 | Predominant Bioaccumulative Toxaphene Congeners (Parlar 26, 50 and 62) | 1.2 mL in nonane |
| Unlabeled (ng/mL) | | |
| Parlar 26 | 2000 | |
| Parlar 50 | 2000 | |
| Parlar 62 | 2000 | |

Multiple Class Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|---|--|------------------|
| NEW ES-5516 | EPA Method 1699 Pesticide Stock Solution | 0.5 mL in nonane |
| Labeled (ng/mL) | | |
| Hexachlorobenzene (¹³ C ₆ , 99%) | 1800 | |
| γ -HCH (γ -BHC) (lindane) (¹³ C ₆ , 99%) | 2600 | |
| Heptachlor (¹³ C ₁₀ , 99%) | 1400 | |
| β -HCH (β -BHC) (¹³ C ₆ , 99%) | 1600 | |
| δ -HCH (δ -BHC) (¹³ C ₆ , 99%) | 1600 | |
| Aldrin (¹³ C ₁₂ , 99%) | 1600 | |
| Oxychlordane (¹³ C ₁₀ , 99%) | 1600 | |
| <i>cis</i> -Heptachlor epoxide (¹³ C ₁₀ , 99%) | 1600 | |
| Endosulfan I (¹³ C ₉ , 99%) | 1600 | |
| Dieldrin (¹³ C ₁₂ , 99%) | 1600 | |
| <i>trans</i> -Chlordane (γ) (¹³ C ₁₀ , 99%) | 1600 | |
| <i>trans</i> -Nonachlor (¹³ C ₁₀ , 98%) | 1600 | |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 1600 | |
| Endrin (¹³ C ₁₂ , 99%) | 1600 | |
| Endosulfan II (¹³ C ₉ , 99%) | 1600 | |
| <i>cis</i> -Nonachlor (¹³ C ₁₀ , 99%) | 1600 | |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 1600 | |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 1600 | |
| Mirex (¹³ C ₁₀ , 99%) | 1600 | |
| Methoxychlor (ring- ¹³ C ₁₂ , 99%) | 1600 | |
| Azinphos-methyl (D ₆ , 98%) | 1600 | |
| Diazinon (diethyl-D ₁₀ , 98%) | 1600 | |
| Fonofos (ring- ¹³ C ₆ , 99%) | 1600 | |
| Atrazine (ring- ¹³ C ₃ , 99%) | 1600 | |
| <i>cis</i> -Permethrin (phenoxy- ¹³ C ₆ , 99%) | 1600 | |
| <i>trans</i> -Permethrin (phenoxy- ¹³ C ₆ , 99%) | 1600 | |
| NEW ES-5560 Multi-Class Pesticide Internal Standard 1.2 mL in acetonitrile | | |
| Labeled (ng/mL) | | |
| Acetochlor (ring- ¹³ C ₆ , 99%) | 1000 | |
| Aldicarb (¹³ C ₂ , 98%; D ₃ , 98%) | 1000 | |
| Chlortoluron (N,N-dimethyl-D ₆ , 98%) | 1000 | |
| Diazinon (diethyl-D ₁₀ , 98%) | 1000 | |
| 2,4-Dichlorophenoxyacetic acid (ring- ¹³ C ₆ , 99%) | 1000 | |
| Simazine (ring- ¹³ C ₃ , 99%) | 1000 | |
| NEW ES-5561 Multi-Class Pesticide Native Standard 1.2 mL in acetonitrile | | |
| Unlabeled (ng/mL) | | |
| Acetochlor | 1000 | |
| Aldicarb | 1000 | |
| Chlortoluron | 1000 | |
| Diazinon | 1000 | |
| 2,4-Dichlorophenoxyacetic acid | 1000 | |
| Simazine | 1000 | |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount | | | | |
|--|---|-----------------------------|-----|-----|-----|-----|
| ES-5348 | POPs Pesticides Calibration Solutions [CS1-CS6] | Set of 6 × 0.2 mL in nonane | | | | |
| <i>Individual calibration solutions are available. Please inquire.</i> | | | | | | |
| Unlabeled | | CS1 | CS2 | CS3 | CS4 | CS5 |
| Hexachlorobenzene | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Aldrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Dieldrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Endrin | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDT | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDE | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 4,4'-DDD | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDT | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDE | 0.4 | 2 | 10 | 40 | 200 | 800 |
| 2,4'-DDD | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Chlordane (γ) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Chlordane (α) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Nonachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Nonachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Oxychlordane | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Heptachlor | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>trans</i> -Heptachlor epoxide | 0.4 | 2 | 10 | 40 | 200 | 800 |
| <i>cis</i> -Heptachlor epoxide | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Mirex | 0.4 | 2 | 10 | 40 | 200 | 800 |
| α -HCH (α -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| β -HCH (β -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| γ -HCH (γ -BHC) (lindane) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| δ -HCH (δ -BHC) | 0.4 | 2 | 10 | 40 | 200 | 800 |
| Cleanup | | | | | | |
| Hexachlorobenzene ($^{13}\text{C}_{6}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Aldrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Dieldrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Endrin ($^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>trans</i> -Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>cis</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Oxychlordane ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Heptachlor ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Mirex ($^{13}\text{C}_{10}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| α -HCH (α -BHC) ($^{13}\text{C}_{6}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| β -HCH (β -BHC) ($^{13}\text{C}_{6}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_{6}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| δ -HCH (δ -BHC) ($^{13}\text{C}_{6}$, 99%) | 20 | 20 | 20 | 20 | 20 | 20 |
| Syringe | | | | | | |
| 4,4'-DiCB ($^{13}\text{C}_{12}$, 99%) (PCB-15) | 20 | 20 | 20 | 20 | 20 | 20 |
| 2,3',4',5-TetraCB ($^{13}\text{C}_{12}$, 99%) (PCB-70) | 20 | 20 | 20 | 20 | 20 | 20 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|------------------------------------|------------------|
| ES-5349 | POPs Pesticides HRMS Cleanup Spike | 1.2 mL in nonane |
| ES-5400 | POPs Cleanup Spike | 1.2 mL in nonane |
| ES-5349-L | POPs Pesticides LRMS Cleanup Spike | 0.5 mL in nonane |

| Labeled | ES-5349 (ng/mL) | ES-5400 (ng/mL) | ES-5349-L (ng/mL) |
|--|--------------------|--------------------|----------------------|
| Hexachlorobenzene (¹³ C ₆ , 99%) | 100 | 200 | 1000 |
| Aldrin (¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| Dieldrin (¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| Endrin (¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 4,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| 2,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 | 200 | 1000 |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| cis-Nonachlor (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| Heptachlor (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| Mirex (¹³ C ₁₀ , 99%) | 100 | 200 | 1000 |
| α -HCH (α -BHC) (¹³ C ₆ , 99%) | 100 | 200 | 1000 |
| β -HCH (β -BHC) (¹³ C ₆ , 99%) | 100 | 200 | 1000 |
| γ -HCH (γ -BHC) (lindane) (¹³ C ₆ , 99%) | 100 | 200 | 1000 |
| δ -HCH (δ -BHC) (¹³ C ₆ , 99%) | 100 | 200 | 1000 |

| | | |
|-----------------|---------------------------------------|------------------|
| ES-5399 | POPs PAR Solution | 1.2 mL in nonane |
| ES-5399-10X-0.5 | POPs PAR Solution (10X concentration) | 0.5 mL in nonane |

| Unlabeled | ES-5399 (ng/mL) | ES-5399-10X-0.5 (ng/mL) |
|--|--------------------|----------------------------|
| Hexachlorobenzene | 200 | 2000 |
| Aldrin | 200 | 2000 |
| Dieldrin | 200 | 2000 |
| Endrin | 200 | 2000 |
| 4,4'-DDT | 200 | 2000 |
| 4,4'-DDE | 200 | 2000 |
| 4,4'-DDD | 200 | 2000 |
| 2,4'-DDT | 200 | 2000 |
| 2,4'-DDE | 200 | 2000 |
| 2,4'-DDD | 200 | 2000 |
| trans-Chlordane (γ) | 200 | 2000 |
| cis-Chlordane (α) | 200 | 2000 |
| trans-Nonachlor | 200 | 2000 |
| cis-Nonachlor | 200 | 2000 |
| Oxychlordane | 200 | 2000 |
| Heptachlor | 200 | 2000 |
| trans-Heptachlor epoxide | 200 | 2000 |
| cis-Heptachlor epoxide | 200 | 2000 |
| Mirex | 200 | 2000 |
| α -HCH (α -BHC) | 200 | 2000 |
| β -HCH (β -BHC) | 200 | 2000 |
| γ -HCH (γ -BHC) (lindane) | 200 | 2000 |
| δ -HCH (δ -BHC) | 200 | 2000 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|----------------------|--|------------------|
| EC-5350 | POPs Pesticides HRMS (PCB) Syringe Spike | 1.2 mL in nonane |
| NEW EC-5350-L | POPs Pesticides HRMS (PCB) Syringe Spike | 1.2 mL in nonane |

| Labeled | EC-5350 (ng/mL) | EC-5350-L (ng/mL) |
|---|--------------------|----------------------|
| 4,4'-DiCB (¹³ C ₁₂ , 99%) (PCB-15) | 100 | 1000 |
| 2,3',4',5-TetraCB (¹³ C ₁₂ , 99%) (PCB-70) | 100 | 1000 |

| | | |
|---------|--|------------------|
| ES-5342 | POPs Pesticides, non-Toxaphene, non-HCH HRMS Cleanup Spike | 1.2 mL in nonane |
|---------|--|------------------|

| Labeled | (ng/mL) |
|---|---------|
| Hexachlorobenzene (¹³ C ₆ , 99%) | 10 |
| Aldrin (¹³ C ₁₂ , 99%) | 100 |
| Dieldrin (¹³ C ₁₂ , 99%) | 20 |
| Endrin (¹³ C ₁₂ , 99%) | 100 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 20 |
| 4,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 |
| 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 20 |
| 2,4'-DDD (ring- ¹³ C ₁₂ , 99%) | 100 |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 10 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 10 |
| cis-Nonachlor (¹³ C ₁₀ , 99%) | 10 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 100 |
| Heptachlor (¹³ C ₁₀ , 99%) | 20 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 20 |
| Mirex (¹³ C ₁₀ , 99%) | 20 |

| | | |
|-----------------|-----------------------------|------------------|
| ES-5344-50X-0.5 | POPs HRMS HCH Cleanup Spike | 0.5 mL in nonane |
|-----------------|-----------------------------|------------------|

| Labeled | (ng/mL) |
|--|---------|
| α -HCH (α -BHC) (¹³ C ₆ , 99%) | 1000 |
| β -HCH (β -BHC) (¹³ C ₆ , 99%) | 1000 |
| γ -HCH (γ -BHC) (lindane) (¹³ C ₆ , 99%) | 1000 |
| δ -HCH (δ -BHC) (¹³ C ₆ , 99%) | 1000 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount | | | | | | | |
|------------------|---|-------------------------------|--|--|--|--|--|--|--|
| ES-5019-A | Persistent Pesticide Calibration Solutions [CS1-CS10] | Set of 10 × 0.25 mL in nonane | | | | | | | |
| ES-5019-A-CS1-8 | Persistent Pesticide Calibration Solutions [CS1-CS8] | Set of 8 × 0.25 mL in nonane | | | | | | | |
| ES-5019-A-CS9-10 | Persistent Pesticide Calibration Solutions [CS9-CS10] | Set of 2 × 0.25 mL in nonane | | | | | | | |

| Unlabeled | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | All concentrations are in ng/mL (ppb) | | | |
|---|-----|-----|-----|-----|-----|-----|---------------------------------------|------|------|------|
| | | | | | | | CS7 | CS8 | CS9 | CS10 |
| Hexachlorobenzene | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| β-HCH (β-BHC) | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| γ-HCH (γ-BHC) (lindane) | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| cis-Heptachlor epoxide (B isomer) | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| Oxychlordane | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| trans-Nonachlor | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| 4,4'-DDE | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | 3000 | 6000 |
| Dieldrin | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| 2,4'-DDT | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | 3000 | 6000 |
| 4,4'-DDT | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| Mirex | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| Dechlorane Plus <i>syn</i> | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| Dechlorane Plus <i>anti</i> | 1.0 | 2.5 | 10 | 35 | 100 | 300 | 500 | 1000 | | |
| Labeled | | | | | | | | | | |
| Hexachlorobenzene (¹³ C ₆ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| Dieldrin (¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| β-HCH (β-BHC) (¹³ C ₆ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| Mirex (¹³ C ₁₀ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 250 | 250 | 250 | 250 | 250 | 250 | 250 | 250 | 250 | 250 |
| Recovery | | | | | | | | | | |
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) (PCB-208) | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) (BDE-77) | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) (BDE-139) | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |

| | | |
|--------------------|---------------------------------------|-----------------------|
| ES-5177-5X10 | Persistent Pesticide Spiking Solution | 5 × 10 mL in methanol |
| ES-5021 | Persistent Pesticide Spiking Solution | 2.5 mL in nonane |
| ES-5177-500X-N-0.5 | Persistent Pesticide Spiking Solution | 0.5 mL in nonane |

| Labeled | ES-5177-5X10 | ES-5021 | ES-5177-500X-N-0.5 |
|---|--------------|---------|--------------------|
| | (ng/mL) | (ng/mL) | (ng/mL) |
| Hexachlorobenzene (¹³ C ₆ , 99%) | 10 | 100 | 5000 |
| Dieldrin (¹³ C ₁₂ , 99%) | 10 | 100 | 5000 |
| β-HCH (β-BHC) (¹³ C ₆ , 99%) | 10 | 100 | 5000 |
| γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%) | 10 | 100 | 5000 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 10 | 100 | 5000 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 10 | 100 | 5000 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 10 | 100 | 5000 |
| Mirex (¹³ C ₁₀ , 99%) | 10 | 100 | 5000 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 10 | 100 | 5000 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 10 | 100 | 5000 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 25 | 250 | 12,500 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount | | | | | | | | |
|-------------|---|-----------------------------|--|--|--|--|--|--|--|--|
| ES-5442 | CDC POPs (with Parlars) Calibration Solutions [CS1-CS9] | Set of 9 x 0.5 mL in nonane | | | | | | | | |

Individual calibration solutions are available. Please inquire.

| Unlabeled | | All concentrations are in ng/mL (ppb) | | | | | | | | |
|--|--|---------------------------------------|-----|-----|-----|-----|-----|------|------|------|
| | | CS1 | CS2 | CS3 | CS4 | CS5 | CS6 | CS7 | CS8 | CS9 |
| Parlar 26 | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Parlar 50 | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Parlar 62 | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Hexachlorobenzene | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| β -HCH (β -BHC) | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| γ -HCH (γ -BHC) (lindane) | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Aldrin | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| cis-Heptachlor epoxide | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Oxychlordane | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| trans-Nonachlor | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| 4,4'-DDE | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | 3000 | 7500 |
| Dieldrin | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Endrin | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Isodrin | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| 2,4'-DDT | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | 3000 | 7500 |
| 4,4'-DDT | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Mirex | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| α -HCH (α -BHC) | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| cis-Chlordanne (α) | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| trans-Chlordanne (γ) | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| 2,4'-DDE | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| cis-Nonachlor | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Methoxychlor | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Pentachloroanisole | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Octachlorostyrene | | 2.5 | 5 | 10 | 30 | 100 | 300 | 1000 | | |
| Labeled | | | | | | | | | | |
| | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Parlar 26 ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Parlar 50 ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Parlar 62 ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Aldrin ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| cis-Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Oxychlordane ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| trans-Nonachlor ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | | 150 | 150 | 150 | 150 | 150 | 150 | 150 | 150 | 150 |
| Dieldrin ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Endrin ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Isodrin ($^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Mirex ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| cis-Chlordanne (α) ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| trans-Chlordanne (γ) ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| cis-Nonachlor ($^{13}\text{C}_{10}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Methoxychlor (ring- $^{13}\text{C}_{12}$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Pentachloroanisole ($^{13}\text{C}_6$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Octachlorostyrene ($^{13}\text{C}_8$, 99%) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| Recovery | | | | | | | | | | |
| | | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
| 1,2,3,4-TetraCDD ($^{13}\text{C}_6$, 99%) | | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 | 25 |
| 2,2',3,3',4,5,5',6,6'-NonaCB ($^{13}\text{C}_{12}$, 99%) (PCB-208) | | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 | 100 |
| 3,3',4,4'-TetraBDE ($^{13}\text{C}_{12}$, 99%) (BDE-77) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |
| 2,2',3,4,4',6-HexaBDE ($^{13}\text{C}_{12}$, 99%) (BDE-139) | | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 | 75 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|-----------------------------|---|------------------|
| NEW ES-5449-10 | CDC POPs (with Parlars) Spiking Standard | 10 mL in nonane |
| NEW ES-5449-100X-1.2 | CDC POPs (with Parlars) Spiking Standard (100X stock) | 1.2 mL in nonane |

| Labeled | ES-5449-10 (ng/mL) | ES-5449-100X-1.2 (ng/mL) |
|---|-----------------------|-----------------------------|
| Parlar 26 (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| Parlar 50 (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| Parlar 62 (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| Hexachlorobenzene (¹³ C ₆ , 99%) | 7.5 | 750 |
| β-HCH (β-BHC) (¹³ C ₆ , 99%) | 7.5 | 750 |
| γ-HCH (γ-BHC) (lindane) (¹³ C ₆ , 99%) | 7.5 | 750 |
| Aldrin (¹³ C ₁₂ , 99%) | 7.5 | 750 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| Oxychlordane (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| trans-Nonachlor (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| 4,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 15.0 | 1500 |
| Dieldrin (¹³ C ₁₂ , 99%) | 7.5 | 750 |
| Endrin (¹³ C ₁₂ , 99%) | 7.5 | 750 |
| Isodrin (¹³ C ₁₂ , 99%) | 7.5 | 750 |
| 2,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 7.5 | 750 |
| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 7.5 | 750 |
| Mirex (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| α-HCH (α-BHC) (¹³ C ₆ , 99%) | 7.5 | 750 |
| cis-Chlordane (α) (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| trans-Chlordane (γ) (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| 2,4'-DDE (ring- ¹³ C ₁₂ , 99%) | 7.5 | 750 |
| cis-Nonachlor (¹³ C ₁₀ , 99%) | 7.5 | 750 |
| Methoxychlor (ring- ¹³ C ₁₂ , 99%) | 7.5 | 750 |
| Pentachloroanisole (¹³ C ₆ , 99%) | 7.5 | 750 |
| Octachlorostyrene (¹³ C ₈ , 99%) | 7.5 | 750 |

| | | |
|-----------------------------|---|--|
| ES-5321 | Multi-Analyte Recovery Spiking Standard | 10 mL in 88% hexane/2% dodecane/10% nonane |
| NEW ES-5321-5X10 | Multi-Analyte Recovery Spiking Standard | 5 × 10 mL in 88% hexane/2% dodecane/10% nonane |
| NEW ES-5321-200X-1.2 | Multi-Analyte Recovery Spiking Standard | 1.2 mL in nonane |

| Labeled | ES-5321 (ng/mL) | ES-5321-200X-1.2 (ng/mL) |
|---|--------------------|-----------------------------|
| 1,2,3,4-TetraCDD (¹³ C ₆ , 99%) | 2.5 | 500 |
| 2,2',3,3',4,5,5',6,6'-NonaCB (¹³ C ₁₂ , 99%) (PCB-208) | 10 | 2000 |
| 3,3',4,4'-TetraBDE (¹³ C ₁₂ , 99%) (BDE-77) | 7.5 | 1500 |
| 2,2',3,4,4',6-HexaBDE (¹³ C ₁₂ , 99%) (BDE-139) | 7.5 | 1500 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|-------------|---|------------------|
| ES-5261-1.2 | Persistent Organic Pollutants Cleanup Spike | 1.2 mL in nonane |

| Labeled | (ng/mL) |
|---|---------|
| Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | 1000 |
| α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | 1000 |
| β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | 1000 |
| γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | 1000 |
| Aldrin ($^{13}\text{C}_{12}$, 99%) | 1000 |
| Dieldrin ($^{13}\text{C}_{12}$, 99%) | 1000 |
| Endrin ($^{13}\text{C}_{12}$, 99%) | 1000 |
| <i>trans</i> -Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | 1000 |
| Oxychlordane ($^{13}\text{C}_{10}$, 99%) | 1000 |
| <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 1000 |
| Heptachlor ($^{13}\text{C}_{10}$, 99%) | 1000 |
| <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | 1000 |
| 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 1000 |
| 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 1000 |
| 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 1000 |

| | | | |
|------------|---------|----------------------------|------------------|
| NEW | ES-5478 | Pesticide Stock Solution 1 | 0.5 mL in nonane |
|------------|---------|----------------------------|------------------|

| Labeled | (ng/mL) |
|---|---------|
| α -HCH (α -BHC) ($^{13}\text{C}_6$, 99%) | 5000 |
| δ -HCH (δ -BHC) ($^{13}\text{C}_6$, 99%) | 5000 |
| Dieldrin ($^{13}\text{C}_{12}$, 99%) | 5000 |
| γ -HCH (γ -BHC) (lindane) ($^{13}\text{C}_6$, 99%) | 5000 |
| <i>cis</i> -Heptachlor epoxide ($^{13}\text{C}_{10}$, 99%) | 5000 |
| Heptachlor ($^{13}\text{C}_{10}$, 99%) | 5000 |
| Oxychlordane ($^{13}\text{C}_{10}$, 99%) | 5000 |
| 2,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| 2,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| 2,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| 4,4'-DDD (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| 4,4'-DDE (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 5000 |
| <i>cis</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 5000 |
| <i>trans</i> -Nonachlor ($^{13}\text{C}_{10}$, 99%) | 5000 |
| <i>trans</i> -Chlordane (γ) ($^{13}\text{C}_{10}$, 99%) | 5000 |
| Mirex ($^{13}\text{C}_{10}$, 99%) | 5000 |

| | | | |
|------------|---------|----------------------------|------------------|
| NEW | ES-5479 | Pesticide Stock Solution 2 | 0.5 mL in nonane |
|------------|---------|----------------------------|------------------|

| Labeled | (ng/mL) |
|--|---------|
| Pentachlorobenzene ($^{13}\text{C}_6$, 99%) | 5000 |
| Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | 5000 |
| β -HCH (β -BHC) ($^{13}\text{C}_6$, 99%) | 5000 |
| Aldrin ($^{13}\text{C}_{12}$, 99%) | 5000 |
| Endosulfan I ($^{13}\text{C}_9$, 99%) | 5000 |
| Endosulfan II ($^{13}\text{C}_9$, 99%) | 5000 |
| Endrin ($^{13}\text{C}_{12}$, 99%) | 5000 |
| <i>cis</i> -Chlordane (α) ($^{13}\text{C}_{10}$, 99%) | 5000 |

Pesticide Standard Mixtures

| Catalog No. | Compound | Amount |
|------------------------|---|------------------|
| NEW ES-5475 | Pesticide Stock Solution 1 | 1.2 mL in nonane |
| | Unlabeled | (ng/mL) |
| | α -HCH (α -BHC) | 5000 |
| | δ -HCH (δ -BHC) | 5000 |
| | Dieldrin | 5000 |
| | γ -HCH (γ -BHC) (lindane) | 5000 |
| | <i>trans</i> -Heptachlor epoxide | 5000 |
| | <i>cis</i> -Heptachlor epoxide | 5000 |
| | Heptachlor | 5000 |
| | Oxychlordane | 5000 |
| | 2,4'-DDD | 5000 |
| | 2,4'-DDE | 5000 |
| | 2,4'-DDT | 5000 |
| | 4,4'-DDD | 5000 |
| | 4,4'-DDE | 5000 |
| | 4,4'-DDT | 5000 |
| | <i>cis</i> -Nonachlor | 5000 |
| | <i>trans</i> -Nonachlor | 5000 |
| | <i>trans</i> -Chlordane (γ) | 5000 |
| | Mirex | 5000 |
| NEW ES-5476 | Pesticide Stock Solution 2 | 1.2 mL in nonane |
| | Unlabeled | (ng/mL) |
| | Pentachlorobenzene | 5000 |
| | Trifluralin | 5000 |
| | Hexachlorobenzene | 5000 |
| | β -HCH (β -BHC) | 5000 |
| | Aldrin | 5000 |
| | Endosulfan I | 5000 |
| | Endosulfan II | 5000 |
| | Endrin | 5000 |
| | <i>cis</i> -Chlordane (α) | 5000 |
| | Endrin aldehyde | 5000 |
| NEW ES-5499-2.2 | PCB/Pollutant Mixture | 2.2 mL in hexane |
| | Labeled | (ng/mL) |
| | 2,4,4'-TriCB ($^{13}\text{C}_{12}$, 99%) (PCB-28) | 10,000 |
| | 2,2',5,5'-TetraCB ($^{13}\text{C}_{12}$, 99%) (PCB-52) | 10,000 |
| | 2,2',4,5,5'-PentaCB ($^{13}\text{C}_{12}$, 99%) (PCB-101) | 10,000 |
| | 2,2',4,4',5,5'-HexaCB ($^{13}\text{C}_{12}$, 99%) (PCB-153) | 10,000 |
| | 4,4'-DDT (ring- $^{13}\text{C}_{12}$, 99%) | 10,000 |
| | Hexachlorobenzene ($^{13}\text{C}_6$, 99%) | 50,000 |
| | 1,2,4,5-Tetrachlorobenzene ($^{13}\text{C}_6$, 99%) | 50,000 |
| NEW ES-5501 | PCB/Pollutant Native Mixture | 1.2 mL in hexane |
| | Unlabeled | (ng/mL) |
| | 2,4,4'-TriCB (PCB-28) | 10,000 |
| | 2,2',5,5'-TetraCB (PCB-52) | 10,000 |
| | 2,2',4,5,5'-PentaCB (PCB-101) | 10,000 |
| | 2,2',4,4',5,5'-HexaCB (PCB-153) | 10,000 |
| | 4,4'-DDT | 10,000 |
| | Hexachlorobenzene | 50,000 |
| | 1,2,4,5-Tetrachlorobenzene | 50,000 |

Pesticide Standard Mixtures

| Catalog No. | Compound | | Amount |
|--------------------|--|-------|------------------|
| NEW ES-5521 | Multi-Functional PAR Spike | | 1.2 mL in nonane |
| | | | |
| | Unlabeled Dioxins | IUPAC | (ng/mL) |
| | 2,3,7,8-TetraCDD | | 200 |
| | 1,2,3,7,8-PentaCDD | | 200 |
| | 1,2,3,4,7,8-HexaCDD | | 200 |
| | 1,2,3,6,7,8-HexaCDD | | 200 |
| | 1,2,3,7,8,9-HexaCDD | | 200 |
| | 1,2,3,4,6,7,8-HeptaCDD | | 200 |
| | OctaCDD | | 200 |
| | 2,3,7,8-TetraCDF | | 200 |
| | 1,2,3,7,8-PentaCDF | | 200 |
| | 2,3,4,7,8-PentaCDF | | 200 |
| | 1,2,3,4,7,8-HexaCDF | | 200 |
| | 1,2,3,6,7,8-HexaCDF | | 200 |
| | 1,2,3,7,8,9-HexaCDF | | 200 |
| | 2,3,4,6,7,8-HexaCDF | | 200 |
| | 1,2,3,4,6,7,8-HeptaCDF | | 200 |
| | 1,2,3,4,7,8,9-HeptaCDF | | 200 |
| | OctaCDF | | 200 |
| | Unlabeled PCBs | | |
| | 2-MonoCB | 1 | 200 |
| | 4-MonoCB | 3 | 200 |
| | 2,2'-DiCB | 4 | 200 |
| | 4,4'-DiCB | 15 | 200 |
| | 2,2',6-TriCB | 19 | 200 |
| | 3,4,4'-TriCB | 37 | 200 |
| | 2,2',6,6'-TetraCB | 54 | 200 |
| | 3,3',4,4'-TetraCB | 77 | 200 |
| | 3,4,4',5-TetraCB | 81 | 200 |
| | 2,2',4,6,6'-PentaCB | 104 | 200 |
| | 2,3,3',4,4'-PentaCB | 105 | 200 |
| | 2,3,4,4',5-PentaCB | 114 | 200 |
| | 2,3',4,4',5-PentaCB | 118 | 200 |
| | 2',3,4,4',5-PentaCB | 123 | 200 |
| | 3,3',4,4',5-PentaCB | 126 | 200 |
| | 2,2',4,4',6,6'-HexaCB | 155 | 200 |
| | 2,3,3',4,4',5-HexaCB | 156 | 200 |
| | 2,3,3',4,4',5'-HexaCB | 157 | 200 |
| | 2,3',4,4',5,5'-HexaCB | 167 | 200 |
| | 3,3',4,4',5,5'-HexaCB | 169 | 200 |
| | 2,2',3,4',5,6,6'-HeptaCB | 188 | 200 |
| | 2,3,3',4,4',5,5'-HeptaCB | 189 | 200 |
| | 2,2',3,3',5,5',6,6'-OctaCB | 202 | 200 |
| | 2,3,3',4,4',5,5',6-OctaCB | 205 | 200 |
| | 2,2',3,3',4,4',5,5',6-NonaCB | 206 | 200 |
| | 2,2',3,3',4,5,5',6,6'-NonaCB | 208 | 200 |
| | DecaCB | 209 | 200 |
| | Unlabeled PBDEs | IUPAC | (ng/mL) |
| | 4-MonoBDE | | 3 200 |
| | 2,4-DiBDE | | 4 200 |
| | 4,4'-DiBDE | | 15 200 |
| | 2,2',4-TriBDE | | 17 200 |
| | 2,4,4'-TriBDE | | 28 200 |
| | 2,2',4,4'-TetraBDE | | 47 200 |
| | 2,2',4,5'-TetraBDE | | 49 200 |
| | 2,3',4,4'-TetraBDE | | 66 200 |
| | 2,3',4',6-TetraBDE | | 71 200 |
| | 3,3',4,4'-TetraBDE | | 77 200 |
| | 2,2',3,4,4'-PentaBDE | | 85 200 |
| | 2,2',4,4',5-PentaBDE | | 99 200 |
| | 2,2',4,4',6-PentaBDE | | 100 200 |
| | 2,3',4,4',6-PentaBDE | | 119 200 |
| | 3,3',4,4',5-PentaBDE | | 126 200 |
| | 2,2',3,4,4',5'-HexaBDE | | 138 400 |
| | 2,2',4,4',5,5'-HexaBDE | | 153 400 |
| | 2,2',4,4',5,6'-HexaBDE | | 154 400 |
| | 2,2',4,4',6,6'-HexaBDE | | 155 400 |
| | 2,3,4,4',5,6-HexaBDE | | 166 400 |
| | 2,2',3,4,4',5,6-HeptaBDE | | 181 400 |
| | 2,2',3,4,4',5,6-HeptaBDE | | 183 400 |
| | 2,3,3',4,4',5,6-HeptaBDE | | 190 400 |
| | 2,2',3,4,4',5,5',6-OctaBDE | | 203 400 |
| | 2,3,3',4,4',5,5',6-OctaBDE | | 205 400 |
| | 2,2',3,3',4,4',5,5',6-NonaBDE | | 206 1000 |
| | 2,2',3,3',4,4',5,6,6'-NonaBDE | | 207 1000 |
| | DecaBDE | | 209 1000 |
| | Unlabeled Pesticides | | |
| | Pentachlorobenzene | | 200 |
| | Hexachlorobenzene | | 200 |
| | α -HCH (α -BHC) | | 200 |
| | β -HCH (β -BHC) | | 200 |
| | γ -HCH (γ -BHC) (lindane) | | 200 |
| | δ -HCH (δ -BHC) | | 200 |
| | Heptachlor | | 200 |
| | <i>cis</i> -Heptachlor epoxide | | 200 |
| | <i>trans</i> -Heptachlor epoxide | | 200 |
| | <i>cis</i> -Nonachlor | | 200 |
| | <i>trans</i> -Nonachlor | | 200 |
| | <i>cis</i> -Chlordane (α) | | 200 |
| | <i>trans</i> -Chlordane (β) | | 200 |
| | Oxychlordane | | 200 |
| | 2,4'-DDD | | 200 |
| | 4,4'-DDD | | 200 |
| | 2,4'-DDE | | 200 |
| | 4,4'-DDE | | 200 |
| | 2,4'-DDT | | 200 |
| | 4,4'-DDT | | 200 |
| | Aldrin | | 200 |
| | Dieldrin | | 200 |
| | Endrin | | 200 |
| | Mirex | | 200 |
| | Chlordecone (kepone) | | 200 |
| | Endosulfan I | | 200 |
| | Endosulfan II | | 200 |

(continued on next page)

Pesticide Standard Mixtures

(continued from previous page)

| NEW | ES-5521 | Multi-Functional PAR Spike (continued) | 1.2 mL in nonane |
|------------------------|---------|--|------------------|
| Unlabeled PAHs | | | |
| | | | (ng/mL) |
| Acenaphthene | | | 200 |
| Acenaphthylene | | | 200 |
| Anthracene | | | 200 |
| Benz[a]anthracene | | | 200 |
| Benzo[b]fluoranthene | | | 200 |
| Benzo[k]fluoranthene | | | 200 |
| Benzol[ghi]perylene | | | 200 |
| Benzo[a]pyrene | | | 200 |
| Chrysene | | | 200 |
| Dibenz[a,h]anthracene | | | 200 |
| Fluoranthene | | | 200 |
| Fluorene | | | 200 |
| Indeno[1,2,3-cd]pyrene | | | 200 |
| Naphthalene | | | 200 |
| Phenanthrene | | | 200 |
| Pyrene | | | 200 |

Chemical Weapon Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|------------------------|---|---|-------------------------------|--------|
| CDNLM-6786-1.2 | Aminomethylphosphonic acid (AMPA) (¹³ C, 99%; ¹⁵ N, 98%, methylene-D ₂ , 98%) | *CH ₄ D ₂ *NO ₃ P | 100 µg/mL in H ₂ O | 1.2 mL |
| CDNLM-6786-10 | | | | 10 mL |
| NEW ULM-9897-1.2 | Diethyl methylphosphonate (unlabeled) | CH ₃ P(O)(OC ₂ H ₅) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9898-1.2 | Diisopropyl methylphosphonate (unlabeled) | C ₇ H ₁₇ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9896-1.2 | Dimethyl methylphosphonate (unlabeled) | C ₃ H ₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW DLM-4851-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9287-M-1.2 | O,O-Diethyl phosphate, potassium salt (DEP) | C ₄ H ₁₀ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9287-M-20X-1.2 | (unlabeled) | | 2 mg/mL in methanol | 1.2 mL |
| DLM-4852-1.2 | O,O-Diethyl thiophosphate, potassium salt (DETP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₃ PS | 100 µg/mL in methanol | 1.2 mL |
| ERD-119 | O,O-Diethyl thiophosphate, potassium salt (DETP) (unlabeled) | C ₄ H ₁₀ KO ₃ PS | 1000 µg/mL in methanol | 1.2 mL |
| NEW DLM-9003-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) (diethyl-D ₁₀ , 98%) | C ₄ D ₁₀ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9002-1.2 | O,O-Diethyl dithiophosphate, potassium salt (DEDTP) | C ₄ H ₁₀ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9002-20X-1.2 | (unlabeled) | | 2 mg/mL in methanol | 1.2 mL |
| NEW DLM-8868-1.2 | O,O-Dimethyl phosphoric acid, potassium salt (DMP) (dimethyl-D ₆ , 98%) | C ₂ D ₆ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8867-1.2 | O,O-Dimethyl phosphoric acid, potassium salt (DMP) | C ₂ H ₆ KO ₄ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8867-20X-1.2 | (unlabeled) | | 2 mg/mL in methanol | 1.2 mL |
| NEW DLM-8904-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) (dimethyl-D ₆ , 98%) CP 97% | C ₂ D ₆ KO ₃ PS | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-8905-1.2 | O,O-Dimethyl thiophosphate, potassium salt (DMTP) | C ₂ H ₆ KO ₃ PS | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-8905-2X-1.2 | (unlabeled) CP 97% | | 2 mg/mL in methanol | 1.2 mL |
| NEW DLM-4541-M-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) (dimethyl-D ₆ , 98%) | C ₂ D ₆ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9004-1.2 | O,O-Dimethyl dithiophosphate, potassium salt (DMDTP) | C ₂ H ₆ KO ₂ PS ₂ | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9004-20X-1.2 | (unlabeled) | | 2 mg/mL in methanol | 1.2 mL |
| NEW ULM-9899-1.2 | Dipinacolyl methylphosphonate (unlabeled) | C ₁₃ H ₂₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9887-1.2 | 1,4-Dithiane (unlabeled) | C ₄ H ₈ S ₂ | 1000 µg/mL in methanol | 1.2 mL |
| ULM-6091-1.2 | Ethyl hydrogen dimethylamidophosphate, sodium salt (unlabeled) CP 90% | C ₄ H ₁₁ NPO ₃ Na | 1000 µg/mL in methanol | 1.2 mL |
| DLM-6098-1.2 | Ethyl hydrogen methylphosphonate (ethyl-D ₅ , 98%) | C ₃ H ₄ D ₅ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-9893-1.2 | N-Ethyldiethanolamine (unlabeled) | (HOCH ₂ CH ₂)NC ₂ H ₅ | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9948-1.2 | Ethylphosphonic acid (unlabeled) | C ₂ H ₅ P(O)(OH) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| ERI-017 | Isopropyl methylphosphonic acid (D ₇ , 98%) | C ₄ H ₄ D ₇ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9892-1.2 | N-Methyldiethanolamine (unlabeled) | CH ₃ N(CH ₂ CH ₂ OH) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| CDLM-6100-1.2 | Methylphosphonic acid (¹³ C, 99%; methyl-D ₃ , 98%) | *CD ₃ H ₂ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| DLM-6196-1.2 | Methylphosphonic acid (methyl-D ₃ , 98%) | CD ₃ H ₂ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ERM-038 | Methylphosphonic acid (unlabeled) | CH ₃ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| CLM-6620-1.2 | Methylphosphonic acid, mono-(1,2,2-trimethylpropyl) ester (trimethylpropyl- ¹³ C ₆ , 99%) | *C ₆ CH ₁₇ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-6619-1.2 | Methylphosphonic acid, mono-(1,2,2-trimethylpropyl) ester (unlabeled) | CH ₃ P(O)(OH)OCH ₂ (CH ₃) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| NEW CLM-6096-1.2 | Methyphosphonic acid, monocyclohexyl ester (cyclohexyl- ¹³ C ₆ , 99%) | C*C ₆ H ₁₅ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-6097-1.2 | Methyphosphonic acid, monocyclohexyl ester (unlabeled) | C ₇ H ₁₅ O ₃ P | 100 µg/mL in methanol | 1.2 mL |
| NEW ULM-6097-10X-1.2 | CP 90% | | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-6099-1.2 | Methylphosphonic acid, monoethyl ester (unlabeled) | CH ₃ P(O)(OH)OCH ₂ CH ₃ | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-6093-1.2 | Methylphosphonic acid, monoisopropyl ester (unlabeled) | C ₄ H ₁₁ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9894-1.2 | Pinacolyl alcohol (unlabeled) | (CH ₃) ₃ CCH(OH)CH ₃ | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9895-1.2 | 1-Propylphosphonic acid (unlabeled) | C ₃ H ₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9886-1.2 | 2-Propylphosphonic acid (unlabeled) | C ₃ H ₉ O ₃ P | 1000 µg/mL in methanol | 1.2 mL |
| CLM-6106-1.2 | Ricinine (ring- ¹³ C ₅ , 99%; cyano- ¹³ C, 99%) | C ₂ *C ₆ H ₈ N ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| NEW ULM-9225-1.2 | Ricinine (unlabeled) | C ₈ H ₈ N ₂ O ₂ | 100 µg/mL in acetonitrile | 1.2 mL |
| ERT-054 | Thiodiglycol (D ₈ , 98%) | C ₄ H ₈ D ₈ O ₂ S | 1000 µg/mL in methanol | 1.2 mL |

Chemical Weapon Metabolite Standards

| Catalog No. | Compound | Formula | Concentration | Amount |
|-------------------------|--|---|------------------------|--------|
| ERT-053 | Thiodiglycol (unlabeled) | C ₄ H ₁₀ O ₂ S | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9890-1.2 | Thiodiglycol sulfone (unlabeled) ~70% WT in H ₂ O | O ₂ S(CH ₂ CH ₂ OH) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| ERT-054 | Thiodiglycol (D ₈ , 98%) | C ₄ H ₁₂ D ₈ O ₂ S | 1000 µg/mL in methanol | 1.2 mL |
| ERT-053 | Thiodiglycol (unlabeled) | C ₄ H ₁₀ O ₂ S | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9890-1.2 | Thiodiglycol sulfone (unlabeled) ~70% WT in H ₂ O | O ₂ S(CH ₂ CH ₂ OH) ₂ | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9889-1.2 | Thiodiglycol sulfoxide (unlabeled) | C ₄ H ₁₀ O ₃ S | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9888-1.2 | 1,4-Thioxane (unlabeled) | C ₄ H ₈ OS | 1000 µg/mL in methanol | 1.2 mL |
| NEW ULM-9891-1.2 | Triethanolamine (unlabeled) | (HOCH ₂ CH ₂) ₃ N | 1000 µg/mL in methanol | 1.2 mL |

Chemical Weapon Metabolite Standard Mixtures

| Catalog No. | Compound | Amount |
|--|--|--------------------|
| NEW ES-5547 | Dialkyl Phosphate and Phosphorothioate Cocktail (D, 98%) | 1.2 mL in methanol |
| Labeled | | |
| <u>O,O-Dimethyl phosphate, potassium salt (dimethyl-D₆, 98%)</u> 10,000 | | |
| <u>O,O-Diethyl phosphate, potassium salt (diethyl-D₁₀, 98%)</u> 10,000 | | |
| <u>O,O-Dimethyl thiophosphate, potassium salt (dimethyl-D₆, 98%)</u> 10,000 | | |
| <u>O,O-Diethyl thiophosphate, potassium salt (diethyl-D₁₀, 98%)</u> 10,000 | | |
| <u>O,O-Dimethyl dithiophosphate, potassium salt (dimethyl-D₆, 98%)</u> 10,000 | | |
| <u>O,O-Diethyl dithiophosphate, potassium salt (diethyl-D₁₀, 98%)</u> 10,000 | | |
| Unlabeled | | |
| <u>O,O-Dimethyl phosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl phosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Dimethyl thiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl thiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Dimethyl dithiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl dithiophosphate, potassium salt</u> 10,000 | | |
| NEW ES-5548 | Dialkyl Phosphate and Phosphorothioate Cocktail | 1.2 mL in methanol |
| Unlabeled | | |
| <u>O,O-Dimethyl phosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl phosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Dimethyl thiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl thiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Dimethyl dithiophosphate, potassium salt</u> 10,000 | | |
| <u>O,O-Diethyl dithiophosphate, potassium salt</u> 10,000 | | |
| NEW ES-5562 | HD Metabolites Native Standards Mixture | 1.2 mL in methanol |
| Unlabeled | | |
| <u>1,4-Dithiane</u> 100 | | |
| <u>Thiodiglycol</u> 100 | | |
| <u>1,4-Thioxane</u> 100 | | |
| <u>Thiodiglycol sulfoxide</u> 100 | | |
| <u>Thiodiglycol sulfone</u> 100 | | |

Chemical Weapon Metabolite Standard Mixtures

| Catalog No. | Compound | Amount |
|--------------------|---|--------------------|
| NEW ES-5563 | Amines and Alcohol Native Standards Mixture | 1.2 mL in methanol |
| | Unlabeled | (μ g/mL) |
| | Triethanolamine | 100 |
| | N-Methyldiethanolamine | 100 |
| | N-Ethyldiethanolamine | 100 |
| | Pinacolyl alcohol | 100 |
| NEW ES-5564 | Phosphonic Acid Native Standards Mixture | 1.2 mL in methanol |
| | Unlabeled | (μ g/mL) |
| | Methylphosphonic acid | 100 |
| | Ethylphosphonic acid | 100 |
| | 1-Propylphosphonic acid | 100 |
| | 2-Propylphosphonic acid | 100 |
| NEW ES-5565 | Phosphonic Acid Esters Native Standards Mixture | 1.2 mL in methanol |
| | Unlabeled | (μ g/mL) |
| | Dimethyl methylphosphonate | 100 |
| | Diethyl methylphosphonate | 100 |
| | Diisopropyl methylphosphonate | 100 |
| | Dipinacolyl methylphosphonate | 100 |
| NEW ES-5566 | Phosphonic Acid Half-Esters Native Standards Mixture | 1.2 mL in methanol |
| | Unlabeled | (μ g/mL) |
| | Methylphosphonic acid, monoethyl ester | 100 |
| | Methylphosphonic acid, monoisopropyl ester | 100 |
| | Methylphosphonic acid, mono-(1,2,2-trimethylpropyl) ester | 100 |
| | Methylphosphonic acid, monocyclohexyl ester | 100 |

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| Acenaphthene (D ₁₀ , 98%) | 149, 152, 153, 179 | Benz[e]aceanthrylene/Benz[j]aceanthrylene (¹³ C ₆ , 94%; D ₈ , 94%) | 147, 151 |
| Acenaphthene (D ₁₀ , 99%) | 153, 200 | Benz[e]aceanthrylene/Benz[j]aceanthrylene (unlabeled) | 150, 154 |
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| Acenaphthylene (¹³ C ₆ , 99%) | 147, 151 | Benzene (¹³ C ₆ , 99%; D ₈ , 98%) | 187 |
| Acenaphthylene (D ₈ , 98%) | 149, 152, 198 | Benzene (D ₈ , 99.5%) | 187, 198, 199, 201 |
| Acenaphthylene (unlabeled) | 150, 154, 155, 237 | Benzidine (ring-D ₈ , 98%) | 187 |
| Acephate (D ₈ , 98%) | 212, 216 | Benzo[a]pyrene (¹³ C ₄ , 99%) | 147, 151, 179 |
| Acephate (unlabeled) | 212, 216 | Benzo[a]pyrene (D ₁₂ , 98%) | 149, 152, 153, 179, 198 |
| Acetaminophen (acetyl- ¹³ C ₂ , 99%; ¹⁵ N, 98%) | 169 | Benzo[a]pyrene (unlabeled) | 150, 154, 155, 179, 237 |
| Acetaminophen (unlabeled) | 169 | (+/-)-Benzo[a]pyrene R-7,T-8,C-9,C-10-tetrahydrotetrol (ring- ¹³ C ₆ , 99%) | 156 |
| Acetamiprid (pyridylmethyl- ¹³ C ₆ , 99%) | 213, 216 | (+/-)-Benzo[a]pyrene R-7,T-8,C-9,T-10-tetrahydrotetrol (ring- ¹³ C ₆ , 99%) | 156 |
| Acetamiprid (unlabeled) | 213, 216 | (+/-)-Benzo[a]pyrene R-7,T-8,T-9,C-10-tetrahydrotetrol (ring- ¹³ C ₆ , 99%) | 156 |
| Acetochlor (ring- ¹³ C ₆ , 99%) | 216, 227 | (+/-)-Benzo[a]pyrene R-7,T-8,T-9,T-10-tetrahydrotetrol (ring- ¹³ C ₆ , 99%) | 156 |
| Acetochlor (unlabeled) | 216, 227 | Benzo[b]fluoranthene (¹³ C ₆ , 99%) | 147, 151, 179 |
| Acetone (D ₈ , 99.9%) | 187 | Benzo[b]fluoranthene (D ₁₂ , 98%) | 149, 152, 153, 179 |
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| Acrylamide (+100 ppm hydroquinone) (1,2,3- ¹³ C ₃ , 99%) | 171 | Benzo[b]furan (unlabeled) | 150 |
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| Acrylonitrile (¹³ C ₃ , 99%) | 187 | Benzo[c]fluorene (unlabeled) | 150 |
| Acrylonitrile (D ₃ , 98%) | 187 | Benzo[c]phenanthrene (¹³ C ₆ , 99%) | 147, 151, 179 |
| Alachlor (ring- ¹³ C ₆ , 99%) | 179, 216 | Benzo[c]phenanthrene (unlabeled) | 150, 154, 179 |
| Aldicarb (¹³ C ₂ , 98%; D ₃ , 98%) | 214, 216, 227 | Benzo[e]pyrene (¹³ C ₄ , 99%) | 147, 179 |
| Aldicarb (unlabeled) | 214, 216, 227 | Benzo[e]pyrene (D ₁₂ , 98%) | 149, 153, 179 |
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| 3-Amino-2-oxazolidone (AOZ) (unlabeled) | 171 | Benzo[j]fluoranthene (unlabeled) | 150, 179 |
| 1-Amino-2-propanol (D ₈ , 98%) | 187 | Benzo[k]fluoranthene (¹³ C ₆ , 99%) | 147, 151, 179 |
| 1-Aminohydantoin hydrochloride (AHD) (5,5-D ₂ , 98%) | 171 | Benzo[k]fluoranthene (D ₁₂ , 98%) | 149, 152, 153, 179, 200 |
| 1-Aminohydantoin hydrochloride (AHD) (unlabeled) | 171 | Benzo[k]fluoranthene (unlabeled) | 150, 154, 155, 237 |
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| 2-Aminonaphthalene (unlabeled) | 187 | Benzyl butyl phthalate (ring-D ₈ , 98%) | 172, 179 |
| Amitriptyline-HCl (N,N-dimethyl-D ₆ , 98%) | 169 | Benzyl butyl phthalate (unlabeled) | 172 |
| Amitriptyline-HCl (unlabeled) | 169 | Benzyl paraben (benzyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | 165, 182 |
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| Ammeline (desethyldeisopropylhydroxyatrazine) (unlabeled) | 171 | Biphenyl (D ₁₀ , 98%) | 187 |
| Amoxicillin-3H ₂ O (phenyl- ¹³ C ₆ , 99%) | 170 | Biphenyl (unlabeled) | 120, 187 |
| Anabasine (2,2',3,4,5,6- ¹³ C ₆ , 99%) | 170 | 1,2-Bis(2,4,6-tribromophenoxy) ethane (¹³ C ₁₂ , 99%) | 142, 143 |
| Anabasine (unlabeled) | 170 | 1,2'-Bis(2,4,6-tribromophenoxy) ethane (unlabeled) | 141, 142 |
| Anatabine (2,2',3,4,5,6- ¹³ C ₆ , 99%) | 170 | Bis(2-chloroethoxy) methane (chloroethoxy-D ₈ , 98%) | 187, 201 |
| Anatabine (unlabeled) | 170 | Bis(2-chloroethyl) ether (D ₈ , 98%) | 187, 198, 200 |
| 4-Androstene-3,17-dione (2,3,4- ¹³ C ₃ , 98%) | 166 | Bis(2-chloroisopropyl) ether (D ₁₂ , 95%) | 187 |
| 4-Androstene-3,17-dione (2,2,4,6-D ₅ , 98%) | 166 | Bis(2-chloroisopropyl) ether (unlabeled) | 187 |
| 4-Androstene-3,17-dione (unlabeled) | 166 | Bis(2-ethylhexyl) adipate (adipate- ¹³ C ₆ , 99%) | 172, 179 |
| Androsterone glucuronide (2,2,4,4-D ₄ , 98%) | 166 | Bis(2-ethylhexyl) adipate (unlabeled) | 172 |
| Androsterone glucuronide (unlabeled) | 166 | Bis(2-ethylhexyl) phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | 172 |
| 1,6-Anhydro- β -D-glucose (levoglucosan) (¹³ C ₆ , 98%) | 171 | Bis(2-ethylhexyl) phthalate (ring-D ₄ , 98%) | 172, 179, 201 |
| Aniline (¹³ C ₆ , 99%) | 187 | Bis(2-ethylhexyl) phthalate (unlabeled) | 172, 179 |
| Aniline (D ₈ , 98%) | 187 | 1,2-Bis(pentabromophenyl) ethane (¹³ C ₁₄ , 99%) | 142, 143 |
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| Anthracene (¹³ C ₆ , 99%) | 147, 151, 179 | 2,4'-Bisphenol A (2-(2-hydroxyphenyl)-2-(4-hydroxyphenyl)propane) (unlabeled) | 141, 142 |
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| Bendiocarb (¹³ C ₃ , 99%) | 214, 216 | Bisphenol F (ring- ¹³ C ₁₂ , 99%) | 174 |
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| 4'-Bromo-2,3,3',4-tetraCB (¹³ C ₁₂ , 99%) | 120, 121 | 5-Chloro-2-(2,4-dichlorophenoxy)-anisole (methyl triclosan) (ring- ¹³ C ₁₂ , 99%) | 202, 204 |
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| Carbaryl (unlabeled) | 179, 214, 216 | 1-Chloropyrene (unlabeled) | 156 |
| Carbazole (ring-D ₉ , 98%) | 187, 200 | 2-Chlorosyringaldehyde (unlabeled) | 197 |
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| 4,4'-DDT (ring- ¹³ C ₁₂ , 99%) | 180, 210, 217, 222, 223, 224, 228, 229, 230, 231, 232, 233, 234 | Dibenzo[furan] (D ₈ , 98%) | 188 |
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| 4,6-Dichloroguaiacol (unlabeled) | 196, 197 | Di-n-butyl phthalate (ring-D ₄ , 98%) | 172, 180, 200 |
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| 1,5-Dichloronaphthalene (DiCN) (¹³ C ₁₀ , 99%) | 160 | Di-n-hexyl phthalate (unlabeled) | 172, 180 |
| 1,5-Dichloronaphthalene (DiCN) (unlabeled) | 160 | 4,6-Dinitro-2-methylphenol (ring-D ₂ , 98%) | 189, 198, 200 |
| 1,8-Dichloronaphthalene (DiCN) (unlabeled) | 160 | 1,3-Dinitrobenzene (¹³ C ₆ , 99%) | 185 |
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| 2,4-Dichlorophenol (¹³ C ₆ , 99%) | 184, 195, 196 | 1,3-Dinitrobenzene (unlabeled) | 185 |
| 2,4-Dichlorophenol (D ₄ , 98%) | 177 | 2,4-Dinitrophenol (ring-D ₃ , 98%) | 185, 189 |
| 2,4-Dichlorophenol (ring-D ₃ , 98%) | 177, 198, 200 | 2,4-Dinitrophenol (unlabeled) | 185, 189 |
| 2,4-Dichlorophenol (unlabeled) | 177, 196 | 2,4-Dinitrotoluene (ring-D ₃ , 98%) | 185, 189 |
| 2,5-Dichlorophenol (¹³ C ₆ , 99%) | 177 | 2,4-Dinitrotoluene (unlabeled) | 185, 189 |
| 2,5-Dichlorophenol (unlabeled) | 177 | 2,6-Dinitrotoluene (methyl-D ₃ , 98%) | 185, 189, 201 |
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| 2,4-Dichlorophenoxyacetic acid (ring-D ₃ , 98%) | 218 | Dinocap (unlabeled) | 218 |
| 2,4-Dichlorophenoxyacetic acid (unlabeled) | 218, 227 | Di-n-octyl phthalate (ring-D ₄ , 98%) | 172, 200 |
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| 1,3-Dichloropropene (D ₄ , 98%) | 189, 198, 199, 201 | Dinotefuran (furylmethyl- ¹³ C ₅ , 99%) | 213, 219 |
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| Dichlorprop (unlabeled) | 218 | Di-n-propyl phthalate (ring-1,2- ¹³ C ₂ , dicarboxyl- ¹³ C ₂ , 99%) | 172, 180 |
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| Endosulfan II (unlabeled) | 180, 209, 210, 219, 222, 223, 225, 235, 236 | Fluorene (D ₁₀ , 98%) | 149, 152, 198, 200 |
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| Glycolithocholic acid, sodium salt (unlabeled) | 167 | 2,2',3,4',5,6-Heptachlorobiphenyl (HeptaCB) (¹³ C ₁₂ , 99%) | 79, 83, 84, 109, 110, 117 |
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| Glyphosate (2- ¹³ C, 99%; ¹⁵ N, 98%+) | 181, 213, 219 | 2,3,3',4,4',5,5'-Heptachlorobiphenyl (HeptaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 101, 109, 111, 112, 114, 115, 118, 120, 236 |
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| α-HCH (α -BHC) (unlabeled) | 210, 216, 222, 223, 225, 228, 229, 232, 235, 236 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran (HeptaCDF) (¹³ C ₁₂ , 99%) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| β-HCH (β -BHC) (¹³ C ₆ , 99%) | 179, 210, 216, 222, 223, 224, 227, 228, 229, 230, 231, 232, 233, 234 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| β-HCH (β -BHC) (unlabeled) | 210, 216, 222, 223, 225, 228, 229, 231, 232, 235, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
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| γ-HCH (γ -BHC) (lindane) (¹³ C ₆ , 99%) | 179, 210, 216, 222, 223, 224, 227, 228, 229, 230, 231, 232, 233, 234 | 1,2,3,4,6,8,9-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 21, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64, 65 |
| γ-HCH (γ -BHC) (lindane) (unlabeled) | 210, 216, 222, 223, 225, 228, 229, 231, 232, 235, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
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| 1,2,3,4,6,7,8-Heptabromodibenzofuran (HeptaBDF) (unlabeled) | 27, 70, 71, 72, 74 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 1,2,3,4,6,7,8-Heptabromodibenzo-p-dioxin (HeptaBDD) (¹³ C ₁₂ , 99%) | 26, 70, 71 | 1,2,3,4,7,8,9-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 1,2,3,4,6,7,8-Heptabromodibenzo-p-dioxin (HeptaBDD) (unlabeled) | 26, 70, 71 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,3',4,5',6-Heptabromodiphenyl ether (HeptaBDE) (unlabeled) | 129 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,4,4',5,6-Heptabromodiphenyl ether (HeptaBDE) (unlabeled) | 129, 132, 133, 136, 138, 139, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,4,4',5,6-Heptabromodiphenyl ether (HeptaBDE) (¹³ C ₁₂ , 99%) | 125, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 143 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,4,4',5'-6-Heptabromodiphenyl ether (HeptaBDE) (unlabeled) | 129, 130, 131, 132, 133, 134, 135, 136, 137, 138, 139, 142, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,3,3',4,4',5,6-Heptabromodiphenyl ether (HeptaBDE) (¹³ C ₁₂ , 99%) | 125 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,3,3',4,4',5,6-Heptabromodiphenyl ether (HeptaBDE) (unlabeled) | 129, 132, 133, 136, 138, 139, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,3,3',4,4',5,6-Heptabromodiphenyl ether (HeptaBDE) (unlabeled) | 181, 209, 211, 219, 222, 223, 224, 227, 228, 229, 230, 234 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| Heptachlor (γ) (unlabeled) | 181, 209, 211, 219, 222, 223, 225, 228, 229, 230, 231, 234 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| cis-Heptachlor epoxide (¹³ C ₁₀ , 99%) | 181, 209, 211, 219, 222, 223, 224, 227, 228, 229, 230, 231, 232, 233, 234 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| cis-Heptachlor epoxide (unlabeled) | 181, 209, 211, 219, 222, 223, 225, 228, 229, 230, 231, 234 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| trans-Heptachlor epoxide (unlabeled) | 209, 211, 219, 222, 223, 225, 228, 229, 235, 236 | 1,2,3,4,6,7,8-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl (HeptaCB) (¹³ C ₁₂ , 99%) | 79, 51, 58, 86, 87, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 102, 103, 113, 115, 116, 117 | 1,2,3,4,6,7,9-Heptachlorodibenzofuran (HeptaCDF) (unlabeled) | 19, 25, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 68, 236 |
| 2,2',3,3',4,4',5-Heptachlorobiphenyl (HeptaCB) (unlabeled) | 81, 50, 89, 90, 94, 96, 97, 98, 101, 109, 111, 112, 114, 115, 116, 118, 119 | 1,2,3,4,5,6,7-Heptachloronaphthalene (HeptaCN) (¹³ C ₁₀ , 99%) | 160, 161, 162 |
| 2,2',3,3',4,4',6-Heptachlorobiphenyl (HeptaCB) (unlabeled) | 81, 118 | 1,2,3,4,5,6,7-Heptachloronaphthalene (HeptaCN) (unlabeled) | 160, 161, 162 |
| 2,2',3,3',4,5,5'-Heptachlorobiphenyl (HeptaCB) (unlabeled) | 81, 112, 114 | 2,2,3,4,5,6,8-Heptachloronaphthalene (HeptaCN) (unlabeled) | 160 |

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| β-Hexabromocyclododecane (unlabeled) | 140 | 2,3,3',4,4',6-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 112, 114, 118 |
| γ-Hexabromocyclododecane (¹³ C ₁₂ , 99%) | 140, 142, 143 | 2,3,3',4,5,5'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 79, 86, 87, 88, 94, 95 |
| γ-Hexabromocyclododecane (unlabeled) | 140, 142 | 2,3,3',4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 119 |
| 1,2,3,4,7,8-Hexabromodibenzofuran (HexaBDF) (¹³ C ₁₂ , 99%) | 27, 70, 71, 72, 73 | 2,3,3',4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 109, 111, 118 |
| 1,2,3,4,7,8-Hexabromodibenzofuran (HexaBDF)(unlabeled) | 27, 70, 71, 72, 74 | 2,3,3',4,4',5,5'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 51, 52, 58, 79, 82, 84, 86, 88, |
| 1,2,3,4,7,8-Hexabromodibenz-p-dioxin (HexaBDD) (¹³ C ₁₂ , 99%) | 26, 70, 71, 72, 73 | 89, 90, 91, 92, 93, 94, 95, 96, | |
| 1,2,3,4,7,8-Hexabromodibenz-p-dioxin (HexaBDD)(unlabeled) | 26, 70, 71, 72, 74 | 97, 98, 99, 100, 109, 110, 113, | |
| 1,2,3,6,7,8-Hexabromodibenz-p-dioxin (HexaBDD) (¹³ C ₁₂ , 99%) | 26, 70, 71, 72, 73, 74 | 116, 117, 158, 159 | |
| 1,2,3,6,7,8-Hexabromodibenz-p-dioxin (HexaBDD)(unlabeled) | 26, 70, 71, 72, 74 | 2,3',4,4',5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, |
| 1,2,3,7,8,9-Hexabromodibenz-p-dioxin (HexaBDD) (¹³ C ₁₂ , 99%) | 26, 70, 71, 72, 73, 74 | 94, 95, 96, 97, 98, 101, 109, 111, | |
| 1,2,3,7,8,9-Hexabromodibenz-p-dioxin (HexaBDD)(unlabeled) | 26, 70, 71, 72, 74 | 112, 114, 115, 118, 119, 236 | |
| 2,2',3,3',4,4'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 128 | 51, 52, 54, 55, 57, 58, 79, 82, | |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether (HexaBDE) (¹³ C ₁₂ , 99%) | 125, 132, 133 | 84, 86, 88, 89, 90, 91, 92, 93, | |
| 2,2',3,4,4',5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 128, 132, 133, 136, 138, 139, 236 | 94, 95, 96, 97, 98, 99, 100, | |
| 2,2',3,4,4',6-Hexabromodiphenyl ether (HexaBDE) | 57, 113, 125, 130, 131, 142, | 109, 110, 183 | |
| (¹³ C ₁₂ , 99%) | 143, 202, 203, 204, 205, 231, | 3,3',4,4',5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 50, 54, 56, 81, 82, 85, 86, 89, |
| 2,2',3,4,4',6-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 232, 233 | 90, 92, 94, 95, 96, 97, 98, 101, | |
| 2,2',3,4,4',6-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 128 | 109, 111, 118, 120, 236 | |
| 2,2',3,4,4',6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 128 | 189 | |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129 | 1,5,5,6,6,10-Hexachlorodecane (¹³ C ₁₀ , 99%) | 175, 181 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (¹³ C ₁₂ , 99%) | 125, 130, 131, 132, 133, 134, | 1,5,5,6,6,10-Hexachlorodecane (unlabeled) | 175, 181 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 135, 136, 137, 138, 142, 143 | 1,2,3,4,6,7-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 130, 131, 132, 133, 134, | 1,2,3,4,6,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 34, 68 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 135, 136, 137, 138, 142, 236 | 1,2,3,4,6,9-Hexachlorodibenzofuran (HexaCDF) (¹³ C ₁₂ , 99%) | 20, 42, 43, 44, 45, 58, 59, 60, 61 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 130, 131, 132, 133, 134, | 1,2,3,4,6,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,4',5,5'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 135, 136, 137, 138, 142, 143 | 1,2,3,4,7,8-Hexachlorodibenzofuran (HexaCDF) (¹³ C ₁₂ , 99%) | 20, 31, 32, 33, 34, 36, 37, 38, |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 130, 131, 132, 133, 134, | 39, 40, 41, 42, 43, 44, 45, 46, | |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 136, 137, 138, 142, 236 | 48, 51, 52, 54, 55, 57, 58, 59, | |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (¹³ C ₁₂ , 99%) | 125, 130, 131, 132, 133, 134, | 60, 61, 62, 63, 64, 65 | |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 135, 136, 137, 138, 142, 143 | 1,2,3,4,7,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 130, 131, 132, 133, 134, | 39, 40, 41, 42, 44, 46, 47, 48, | |
| 2,2',3,4,4',5,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 136, 137, 138, 142, 143 | 49, 50, 54, 56, 59, 60, 61, 62, | |
| 2,2',3,4,4',6,6'-Hexabromodiphenyl ether (HexaBDE) (¹³ C ₁₂ , 99%) | 125 | 66, 67, 236 | |
| 2,2',3,4,4',6,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 132, 133, 136, 138, 139, 236 | 1,2,3,4,7,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,4',6,6'-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 132, 133, 136, 138, 139, 236 | 1,2,3,4,8,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 34, 68 |
| 2,3,4,4',5,6-Hexabromodiphenyl ether (HexaBDE) (unlabeled) | 129, 132, 133, 136, 138, 139, 236 | 1,2,3,6,7,8-Hexachlorodibenzofuran (HexaCDF) (¹³ C ₁₂ , 99%) | 20, 31, 32, 33, 34, 36, 37, 40, |
| Hexachloro-1,3-butadiene (¹³ C ₄ , 99%) | 189 | 41, 42, 43, 44, 45, 46, 51, 52, | |
| Hexachloro-1,3-butadiene (unlabeled) | 189 | 54, 55, 57, 58, 59, 60, 61, 62, | |
| Hexachlorobenzene (¹³ C ₆ , 99%) | 177, 181, 195, 200, 211, 220, | 63, 64, 65 | |
| Hexachlorobenzene (unlabeled) | 222, 223, 224, 227, 228, 229, | 1,2,3,6,7,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, |
| 2,2',3,3',4,4'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 78, 94, 95, 102, 103, 112, | 39, 40, 41, 42, 44, 46, 47, 48, | |
| 2,2',3,3',4,4'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 113, 116 | 49, 50, 54, 56, 59, 60, 61, 62, | |
| 2,2',3,3',4,6'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 112, 114, 118 | 66, 67, 236 | |
| 2,2',3,3',4,6'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 81, 118 | 1,2,3,6,7,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 78, 51, 53, 83, 84, 87, 88, 89, | 1,2,3,6,8,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 90, 91, 92, 93, 96, 97, 98, 99, | 1,2,3,7,8,9-Hexachlorodibenzofuran (HexaCDF) (¹³ C ₁₂ , 99%) | 20, 31, 32, 33, 34, 36, 37, 40, |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 102, 103, 104, 105, 106, 109, 110, | 41, 42, 43, 44, 45, 46, 51, 52, | |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 112, 113, 115, 116, 117, 130, 131 | 54, 55, 57, 58, 59, 60, 61, 62, | |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 87, 102, 103, 104, 105, | 63, 64, 65 | |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 107, 108, 111, 113, 115, 117, 158, | 1,2,3,7,8,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, |
| 2,2',3,4,4',5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 111, 112, 114, 115, 116, 118 | 39, 40, 41, 42, 44, 46, 47, 48, | |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 79, 104, 105, 117 | 49, 50, 54, 56, 59, 60, 61, 62, | |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 119 | 66, 67, 236 | |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 112, 114, 115, 116 | 1,2,4,6,7,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 104, 105, 108, 111, 112, | 1,2,4,6,7,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 114, 118 | 1,2,4,6,8,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 81, 112, 114, 118 | 1,3,4,6,7,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 79, 82, 84, 109, 110, 116 | 1,3,4,6,7,9-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 24 |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 81, 82, 85, 108, 111, 115, 118, | 2,3,4,6,7,8-Hexachlorodibenzofuran (HexaCDF) (¹³ C ₁₂ , 99%) | 20, 31, 32, 33, 34, 36, 37, 40, |
| 2,2',3,4,5,5'-Hexachlorobiphenyl (HexaCB) (unlabeled) | 119, 120, 236 | 43, 44, 45, 46, 51, 52, 54, 55, | |
| 2,2',3,4,4',5-Hexachlorobiphenyl (HexaCB) (¹³ C ₁₂ , 99%) | 51, 52, 58, 79, 82, 84, 86, 88, | 57, 58, 59, 60, 61, 62, 63, 64, 65 | |
| 2,2',3,4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 89, 90, 91, 92, 93, 94, 95, 96, | 2,3,4,6,7,8-Hexachlorodibenzofuran (HexaCDF) (unlabeled) | 23, 24, 31, 32, 33, 34, 36, 38, |
| 2,2',3,4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 97, 98, 99, 100, 109, 110, 113, | 39, 40, 41, 42, 44, 46, 47, 48, | |
| 2,2',3,4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 115, 116 | 49, 50, 54, 56, 59, 60, 61, 62, | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, | 66, 67, 236 | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 94, 95, 96, 97, 98, 101, 108, | 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 51, 52, 62, 63 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 112, 114, 115, 116, 118, 236 | 1,2,3,4,6,7-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 19, 25, 68 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 51, 52, 58, 79, 82, 84, 86, 88, | 1,2,3,4,6,8-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 89, 90, 91, 92, 93, 94, 95, 96, | 1,2,3,4,6,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 97, 98, 99, 100, 109, 110, 113, | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 40, |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 111, 112, 114, 118, 236 | 41, 42, 43, 44, 45, 46, 51, 52, | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, | 54, 55, 57, 58, 59, 60, 61, 62, | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 94, 95, 96, 97, 98, 101, 108, | 63, 64, 65 | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 97, 98, 99, 100, 109, 110, 113, | 1,2,3,4,7,8-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 19, 25, 31, 32, 33, 34, 36, 38, |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, | 39, 40, 41, 42, 44, 46, 47, 48, | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 94, 95, 96, 97, 98, 101, 108, | 49, 50, 54, 56, 59, 60, 61, 62, | |
| 2,3,3',4,4',5-Hexachlorobiphenyl (HexaCB) (unlabeled) | 111, 112, 114, 118, 236 | 66, 67, 236 | |

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| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64, 65 | 2-Hydroxyfluorene (random- ¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,3,6,7,8-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 19, 25, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 236 | 2-Hydroxyfluorene (unlabeled) | 157, 158, 159 |
| 1,2,3,6,7,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 3-Hydroxyfluorene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,3,6,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 3-Hydroxyfluorene (unlabeled) | 157, 158, 159 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64 | 9-Hydroxyfluorene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 19, 25, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 236 | 9-Hydroxyfluorene (unlabeled) | 157, 158, 159 |
| 1,2,3,6,7,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64 | 3-Hydroxymethyltetrahydrofuran (¹³ C ₅ , 99%) | 190 |
| 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 1-Hydroxynaphthalene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,4,6,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 1-Hydroxynaphthalene (unlabeled) | 157, 158, 159 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64 | 2-Hydroxynaphthalene (unlabeled) | 157, 158, 159 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 19, 25, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 60, 61, 62, 66, 67, 236 | 1-Hydroxynaphthalene (¹³ C ₄ , 99%) | 157, 158, 159 |
| 1,2,3,7,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 47, 48, 49, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64 | 1-Hydroxynaphthalene (unlabeled) | 157, 158, 159 |
| 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin / 1,2,4,6,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) isomer pair (unlabeled) | 19, 34, 68 | 2-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,4,6,7,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 2-Hydroxyphenanthrene (unlabeled) | 157, 158, 159 |
| 1,2,4,6,8,9-Hexachlorodibenzo-p-dioxin (HexaCDD) (unlabeled) | 25 | 3-Hydroxyphenanthrene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 2,2',3,4,5,5'-Hexachlorodiphenyl ether (HexaCDE) (¹³ C ₁₂ , 99%) | 121 | 3-Hydroxyphenanthrene (unlabeled) | 157, 158, 159 |
| Hexachloroethane (1- ¹³ C, 99%) | 189, 200 | 4-Hydroxyphenanthrene (¹³ C ₄ , 99%) | 157, 158, 159 |
| 1,2,3,4,5,6-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | 4-Hydroxyphenanthrene (unlabeled) | 157, 158, 159 |
| 1,2,3,4,5,7-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160, 161 | 17 α -Hydroxypregnolone (20,21- ¹³ C ₂ , 99%; 16,16-D ₂ , 99%) | 167 |
| 1,2,3,4,5,7-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161 | 17 α -Hydroxypregnolone (unlabeled) | 167 |
| 1,2,3,4,5,8-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160 | 17 α -Hydroxyprogesterone (2,3,4- ¹³ C ₃ , 98%) | 167 |
| 1,2,3,4,5,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | 17 α -Hydroxyprogesterone (2,2,4,6,6,21,21,21-D ₈ , 98%) | 168 |
| 1,2,3,4,6,7-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161, 162 | 17 α -Hydroxyprogesterone (unlabeled) | 168 |
| 1,2,3,5,6,7-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160, 161, 162 | 1-Hydroxypyrene (¹³ C ₆ , 99%) | 157, 158, 159 |
| 1,2,3,5,6,7-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161, 162 | 1-Hydroxypyrene (unlabeled) | 157, 158, 159 |
| 1,2,3,5,6,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | I | |
| 1,2,3,5,7,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | Ibuprofen (propionic- ¹³ C ₃ , 99%) | 169 |
| 1,2,3,6,7,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | Ibuprofen (unlabeled) | 169 |
| 1,2,3,4,5,6-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160, 161 | Imidacloprid (4,4,5,5-D ₄ , 98%) | 213, 220 |
| 1,2,3,4,5,6-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161 | Imidacloprid (unlabeled) | 213, 220 |
| 1,2,3,4,5,8-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160 | Imipramine-HCl (2,4,6,8-D ₄ , 98%) | 169 |
| 1,2,3,4,5,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | Imipramine (unlabeled) | 169 |
| 1,2,3,4,6,7-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161, 162 | Indeno[1,2,3- <i>c</i>]pyrene (¹³ C ₆ , 99%) | 148, 151, 181 |
| 1,2,3,5,6,7-Hexachloronaphthalene (HexaCN) (¹³ C ₁₀ , 99%) | 160, 161, 162 | Indeno[1,2,3- <i>c</i>]pyrene (D ₁₂ , 98%) | 149, 152, 153, 181 |
| 1,2,3,5,6,7-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161 | Indeno[1,2,3- <i>c</i>]pyrene (unlabeled) | 150, 154, 155, 237 |
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| 1,2,4,5,6,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160, 161, 162 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | 165, 182 |
| 1,2,4,5,7,8-Hexachloronaphthalene (HexaCN) (unlabeled) | 160 | Isobutyl paraben (isobutyl 4-hydroxybenzoate) (unlabeled) | 165, 182 |
| Hexachlorophenol (¹³ C ₁₃ , 99%) | 165, 184 | Isodrin (¹³ C ₁₂ , 99%) | 181, 209, 211, 220, 222, 223, 224, 232, 233 |
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| Hexamethylcyclotrisiloxane "D ₃ " (hexamethyl- ¹³ C ₆ , 98%) | 165 | Isopropyl methylphosphonic acid (D ₇ , 98%) | 238 |
| Hexamethylcyclotrisiloxane "D ₃ " (unlabeled) | 165 | Isopropyl methylphosphonic acid (unlabeled) | 213, 220 |
| n-Hexane (D ₁₄ , 98%) | 186 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (ring- ¹³ C ₆ , 99%) | 165, 182 |
| Hexanoic acid (D ₁₁ , 98%) | 189 | Isopropyl paraben (isopropyl 4-hydroxybenzoate) (unlabeled) | 165, 182 |
| 2-Hexanone (1,1,1,3-D ₅ , 98%) | 198, 199 | J | |
| n-Hexatriacontane (D ₇₄ , 98%) | 186 | JECS Labeled Mixture Solution | 206 |
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| Nitroglycerin (unlabeled) | 185 | 2,2',3,4,4',5,5',6-Octabromodiphenyl ether (OctaBDE) (unlabeled) | 129, 132, 133, 142, 236 |
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| PAH-SIM Recovery Standard Mixture | 153, 155 | 1,2,3,7,8-Pentabromodibenz-p-dioxin (PentaBDD) (¹³ C ₁₂ , 99%) | 26, 70, 71, 72, 73 |
| PAH Surrogate Cocktail | 152 | 1,2,3,7,8-Pentabromodibenz-p-dioxin (PentaBDD) (unlabeled) | 26, 70, 71, 72, 74 |
| PAH Surrogate Standard Mixture | 152 | 2,2',3,4,4'-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 132, 133, 136, 138, 139, 142, 236 |
| PAH Two-Component Mixture | 153 | 2,2',3,4,4'-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 143 |
| Parathion (diethyl-D ₁₀ , 98%) | 182, 213, 220 | 2,2',4,4',5-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 132, 133, 136, 138, 139, 142, 236 |
| Parathion (unlabeled) | 182, 213, 220 | 2,2',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| Parlar 26 (¹³ C ₁₀ , 99%) | 182, 215, 220, 226, 232, 233 | 2,2',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 143 |
| Parlar 26 (unlabeled) | 182, 215, 220, 226, 232 | 2,2',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Cocktail for CARB Method 429 | 152, 154 | 2,2',3,4,4'-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Injection Standard | 153 | 2,3',4,4',5-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Native Standard Mixture | 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Recovery Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH-SIM Recovery Standard Mixture | 153, 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Surrogate Cocktail | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Surrogate Standard Mixture | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| PAH Two-Component Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| Parathion (diethyl-D ₁₀ , 98%) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| Parathion (unlabeled) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (unlabeled) | 128, 130, 131, 132, 133, 134, 135, 136, 137, 138, 142, 236 |
| Parlar 26 (¹³ C ₁₀ , 99%) | 182, 215, 220, 226, 232, 233 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parlar 26 (unlabeled) | 182, 215, 220, 226, 232 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Cocktail for CARB Method 429 | 152, 154 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Injection Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Native Standard Mixture | 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Recovery Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH-SIM Recovery Standard Mixture | 153, 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Cocktail | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Standard Mixture | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Two-Component Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parathion (diethyl-D ₁₀ , 98%) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parathion (unlabeled) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parlar 26 (¹³ C ₁₀ , 99%) | 182, 215, 220, 226, 232, 233 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parlar 26 (unlabeled) | 182, 215, 220, 226, 232 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Cocktail for CARB Method 429 | 152, 154 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Injection Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Native Standard Mixture | 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Recovery Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH-SIM Recovery Standard Mixture | 153, 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Cocktail | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Standard Mixture | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Two-Component Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parathion (diethyl-D ₁₀ , 98%) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parathion (unlabeled) | 182, 213, 220 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parlar 26 (¹³ C ₁₀ , 99%) | 182, 215, 220, 226, 232, 233 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| Parlar 26 (unlabeled) | 182, 215, 220, 226, 232 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Cocktail for CARB Method 429 | 152, 154 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Injection Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Native Standard Mixture | 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Recovery Standard | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH-SIM Recovery Standard Mixture | 153, 155 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Cocktail | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Surrogate Standard Mixture | 152 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, 137, 138, 142, 236 |
| PAH Two-Component Mixture | 153 | 2,3',4,4',6-Pentabromodiphenyl ether (PentaBDE) (¹³ C ₁₂ , 99%) | 125, 136, |

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|--|--|---|--|
| 2,2',3,4,4'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80 | 1,2,6,7,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',3,4,5'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 108, 111, 112, 114, 115, 116, 118 | 1,2,6,7,9-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',3',4,5-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 78 | 1,3,4,6,7-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',3',4,5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80 | 1,3,4,6,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 22, 24, 34, 68 |
| 2,2',3,5',6-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 108, 111, 118 | 1,3,4,6,9-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 108, 111, 112, 114, 115, 116, 118 | 1,3,4,7,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',4,5,5'-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 78, 83, 84, 87, 88, 102, 103, 104, 105, 106, 109, 110, 112, 113, 115, 116, 117, 235 | 1,3,4,7,9-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',4,5,5'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 87, 102, 103, 104, 105, 107, 108, 111, 112, 114, 115, 116, 118, 119, 235 | 1,3,6,7,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',4,6,6'-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 78, 82, 84, 109, 110 | 1,4,6,7,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,2',4,6,6'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 82, 85, 108, 111, 118, 120, 236 | 2,3,4,6,7-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,3,3',4,4'-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 109, 110, 112, 113, 115, 116, 117, 158, 159 | 2,3,4,6,8-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,3,3',4,4'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 80, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 108, 111, 112, 114, 115, 116, 118, 119, 236 | 2,3,4,6,9-Pentachlorodibenzofuran (PentaCDF) (unlabeled) | 24 |
| 2,3,3',4',6-Pentachlorobiphenyl (PentaCB) (unlabeled) | 80, 104, 105, 108, 111, 112, 114, 115, 116, 118 | 1,2,3,4,6-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3,3',5,5'-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 51, 53, 78, 83, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 109, 110, 116, 117 | 1,2,3,4,7-Pentachlorodibenzo-p-dioxin (PentaCDD) (¹³ C ₁₂ , 99%) | 17, 42, 43, 44, 45 |
| 2,3,3',5,5'-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81, 119 | 1,2,3,4,7-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3,3',5,6-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81 | 1,2,3,6,7-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3,4,4',5-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 109, 110, 112, 113, 117 | 1,2,3,6,8-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3,4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 102, 103, 108, 111, 112, 118, 236 | 1,2,3,6,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 109, 110, 112, 113, 117 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin (PentaCDD) (¹³ C ₁₂ , 99%) | 17, 31, 32, 33, 34, 36, 37, 38, 39, 40, 41, 42, 43, 44, 45, 46, 48, 51, 52, 54, 55, 57, 58, 59, 60, 61, 62, 63, 64, 65 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (¹³ C ₁₂ , 99%) | 81, 119 | 1,2,3,7,8-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 18, 25, 31, 32, 33, 34, 36, 38, 39, 40, 41, 42, 44, 46, 47, 48, 49, 50, 54, 56, 59, 62, 66, 67, 236 |
| 2,3,4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 102, 103, 108, 111, 112, 118, 236 | 1,2,3,7,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 104, 105, 106, 109, 110, 112, 113, 115, 116, 117 | 1,2,3,8,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 18, 25, 34, 68 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81 | 1,2,4,6,7-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 109, 110, 112, 113, 117 | 1,2,4,6,8-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81 | 1,2,4,6,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 19, 34, 68 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 102, 103, 108, 111, 112, 118, 236 | 1,2,4,7,8-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 104, 105, 106, 109, 110, 112, 113, 115, 116, 117 | 1,2,4,7,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81 | 1,2,4,8,9-Pentachlorodibenzo-p-dioxin (PentaCDD) (unlabeled) | 25 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 111, 112, 118, 236 | 1,2,3,4,5-Pentachloronaphthalene (PentaCN) (unlabeled) | 160 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 104, 105, 106, 109, 110, 112, 113, 115, 116, 118, 236 | 1,2,3,4,6-Pentachloronaphthalene (PentaCN) (unlabeled) | 160 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 109 | 1,2,3,5,7-Pentachloronaphthalene (PentaCN) (¹³ C ₁₀ , 99%) | 160, 161, 162 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 108 | 1,2,3,5,7-Pentachloronaphthalene (PentaCN) (unlabeled) | 160, 161, 162 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 110, 112, 113 | 1,2,3,5,8-Pentachloronaphthalene (PentaCN) (unlabeled) | 160, 162 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 102, 103, 104, 105, 107, 108, 111, 112, 114, 115, 116, 118, 236 | 1,2,3,6,7-Pentachloronaphthalene (PentaCN) (unlabeled) | 160 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 54, 55, 57, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 109, 110, 112, 118, 236 | 1,2,4,5,8-Pentachloronaphthalene (PentaCN) (unlabeled) | 160 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 54, 55, 57, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 109, 110, 112, 118, 236 | 1,2,4,6,7-Pentachloronaphthalene (PentaCN) (unlabeled) | 160 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 109 | Pentachloronitrobenzene (¹³ C ₆ , 99%) | 177, 183 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 108 | Pentachloronitrobenzene (unlabeled) | 177, 183 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 51, 52, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 102, 103, 104, 105, 106, 109, 110, 112, 113, 115, 116, 117 | Pentachlorophenol (¹³ C ₆ , 99%) | 177, 183, 195, 196, 200, 203, 205 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 81 | Pentachlorophenol (unlabeled) | 177, 183, 196 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 102, 103, 108, 111, 112, 118, 236 | n-Pentadecane (D ₃₂ , 98%) | 186 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 94, 95, 96, 97, 98, 99, 100, 109, 110, 183 | n-Pentane (D ₁₂ , 98%) | 186 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 54, 56, 81, 82, 85, 86, 89, 90, 92, 94, 95, 96, 97, 98, 99, 101, 108, 111, 118, 120, 236 | Perchloric acid, sodium salt (¹⁸ O ₄ , 90%+) | 171, 190 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 94, 95, 96, 97, 98, 99, 100, 109, 110, 183 | Perchloric acid, sodium salt (unlabeled) | 171, 190 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 50, 51, 52, 59, 62, 63 | Perfluorobutyric acid (PFBA) (unlabeled) | 174 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 24 | Perfluorodecanoic acid (PFDA) (¹³ C ₉ , 99%) | 174 |
| 2,3',4,4',5-Pentachlorobiphenyl (PentaCB) (unlabeled) | 24 | Perfluorodecanoic acid (PFDA) (unlabeled) | 174 |
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| 2,3,4',5-Tetrachlorobiphenyl (TetraCB) (¹³ C ₁₂ , 99%) | 51, 53, 58, 78, 89, 90, 91, 92, 93, 93, 96, 97, 98, 99 | 1,2,4,6-Tetrachlorodibenzo-p-dioxin (TetraCDD) (unlabeled) | 25 |
| 2,3,4',5-Tetrachlorobiphenyl (TetraCB) (unlabeled) | 80, 108, 111, 118 | 1,2,4,7-Tetrachlorodibenzo-p-dioxin (TetraCDD) (unlabeled) | 25 |
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| 2,3,4',5-Tetrachlorobiphenyl (TetraCB) (unlabeled) | 80 | 1,2,4,9-Tetrachlorodibenzo-p-dioxin (TetraCDD) (unlabeled) | 25 |
| 2,3,4',5-Tetrachlorobiphenyl (TetraCB) (unlabeled) | 51, 52, 54, 55, 57, 58, 78, 82, 84, 86, 88, 89, 90, 91, 92, 93, 94, 95, 96, 97, 98, 99, 100, 109, 110, 111, 118, 236 | 1,2,6,7-Tetrachlorodibenzo-p-dioxin (TetraCDD) (unlabeled) | 25 |
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