

3855 NORTH OCOEE STREET SUITE 200, CLEVELAND, TN 37312 OFFICE: (423) 336-4000 FAX: (423) 336-4166

November 10, 2009

Mr. Michael J. Hinton, P.E. Environmental Engineer II New York State Department of Environmental Conservation 270 Michigan Avenue Buffalo, New York 14203-2999

Subject: Charles Gibson Site NYSDEC Registry No. 9-32-063 Stream Sediment Data

RECEIVED

NOV 1 6 2009

NYSDEC REG 9 FOIL UNREL

Dear Mr. Hinton:

Olin hereby submits the annual 2009 Stream Sediment data for the Gibson site in Niagara Falls, NY. These data were collected in September, 2009, and are being submitted ahead of the annual report per your request, due to elevated pesticide levels detected in the annual 2008 sample.

As the data show, both the upstream and downstream pesticide levels have shown significant decreases relative to 2008 levels. The 2009 data show upstream sediment levels to be equal or greater than down stream levels. The attached CD contains the original laboratory report and the data validation memo. A tabulated data set for Gibson site sediment is attached.

Sincerely,

Michael Billoth

OLIN CORPORATION Michael J. Bellotti Principal Environmental Specialist

cc: Curt Richards - Olin via e-mail Matthew Forcucci – NYSDOH Buffalo

Olin Gibson Site Stream Sediment sampling data 2007, 08, 09

Olin Gibson Site, Niagara	Falls, NY	Sediment sampl	es	Sep-07							
Job No	Client Sample ID	Lab Smp ID	Samp Date	Recvd Date	Anal Date	CAS No		Parameter N	Flags	Result	UM
upstream dup	MS-1-091307	A7A33707	9/13/2007	9/13/2007	9/19/2007	319-84-6	alpha-BHC			40	UG/KG
upstream dup	MS-1-091307	A7A33707	9/13/2007	9/13/2007	9/19/2007	319-85-7	beta-BHC		J	4.8	UG/KG
upstream dup	MS-1-091307	A7A33707	9/13/2007	9/13/2007	9/19/2007	319-86-8	delta-BHC		J	4.6	UG/KG
upstream dup	MS-1-091307	A7A33707	9/13/2007	9/13/2007	9/19/2007	58-89-9	gamma-BHC	(Lindane)	J	3.7	UG/KG
							-				
Upstream	US-1-091307	A7A33706	9/13/2007	9/13/2007	9/19/2007	319-84-6	alpha-BHC			39	UG/KG
	US-1-091307	A7A33706	9/13/2007	9/13/2007	9/19/2007	319-85-7	beta-BHC			8	UG/KG
	US-1-091307	A7A33706	9/13/2007	9/13/2007	9/19/2007	319-86-8	delta-BHC			55	UG/KG
	US-1-091307	A7A33706	9/13/2007	9/13/2007	9/19/2007	58-89-9	gamma-BHC	(Lindane)	.1	3.5	UG/KG
							3	()		0.0	00/110
Downstream											
Olin Gibson Site, Niagara	Falls, NY	Sediment sampl	es	Sep-08							
			No.					Parameter		Constant of the	
Job No	Client Sample ID	Lab Smp ID	Samp Date	Recyd Date	Anal Date	CAS No	Analyte	Name	Flags	Result	Unite
Upstream	US-1-091108	A8B13008	9/11/2008	9/11/2008	9/29/2008	319-86-8	delta-BHC	Hume	lingo	17	LIG/KG
-	US-1-091108	A8B13008	9/11/2008	9/11/2008	9/29/2008	58-89-9	damma_BHC	(Lindana)	5	26	UG/KG
	US-1-091108	A8B13008	9/11/2008	9/11/2008	9/29/2009	310 84 6	alaba BUC	(Lindane)	0	20	UG/KG
	US-1-091108	A8B13008	9/11/2008	9/11/2008	9/29/2008	210 95 7	hoto BHC			11	UG/KG
	03-1-031100	A0B13000	9/11/2008	9/11/2006	9/29/2008	319-05-7	Deta-BHC			69	UG/KG
upstream duo	MS 1 001109	A9P12000	0/11/2000	0/11/2000	0/20/2000	240.00.0	date DUO				
upstream dup	MS 1 001109	A0D13009	9/11/2008	9/11/2008	9/29/2008	519-00-0			J	25	UG/KG
	MG-1-091108	A0D13009	9/11/2008	9/11/2008	9/29/2008	58-89-9	gamma-BHC	(Lindane)	U	52	UG/KG
upstream dup	MS-1-091108	A6B13009	9/11/2008	9/11/2008	9/29/2008	319-85-7	beta-BHC			89	UG/KG
upstream dup	MS-1-091108	A8B13009	9/11/2008	9/11/2008	9/29/2008	319-84-6	alpha-BHC			82	UG/KG
Description	D0 1 001100	10010010									
Downstream	DS-1-091108	A8B13010	9/11/2008	9/11/2008	9/29/2008	58-89-9	gamma-BHC	(Lindane)	J	82	UG/KG
	DS-1-091108	A8B13010	9/11/2008	9/11/2008	9/29/2008	319-86-8	delta-BHC		J	66	UG/KG
	DS-1-091108	A8B13010	9/11/2008	9/11/2008	9/29/2008	319-84-6	alpha-BHC			5200	UG/KG
	DS-1-091108	A8B13010	9/11/2008	9/11/2008	9/29/2008	319-85-7	beta-BHC			1000	UG/KG
Olin Gibson Site, Niagara	Falls, NY	Sediment sample	es	Sep-09							
						CAS					
	SAMPLENAME	LABSAMPID	SAMP DATE	PREPDATE	ANADATE	NUMBER	ANALYTE	1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 - 1990 -	Flags	Result	UNITS
Linetre em	000 1104 004700	D010004.04	00/17/0000								
opstream	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC			240	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC [2C]			270	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-85-7	beta-BHC [2C]			280	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-85-7	beta-BHC			260	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-86-8	delta-BHC	1	J	39	ua/ka
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	319-86-8	delta-BHC [2C]		ľ.	42	
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2000	59 90 0	gamma BHC (Lindana)		5	42	ug/kg
	CGS-US1-091709	PSI0634 01	09/17/2009	00/21/2009 00	09/29/200	50-09-9	gamma-BHC (Lindane)		J	10	ug/kg
	CCS US1 001709	RS10034-01	09/17/2009	09/21/2009 08	09/29/200	56-69-9	gamma-BHC (Lindane) [2C]		J	26	ug/kg
	CG3-031-091709	RS10634-01	09/17/2009	09/21/2009 08	09/29/200	2051-24-3	Decachlorobiphenyl			41.3	ug/kg
	CGS-051-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	2051-24-3	Decachlorobiphenyl [2C]			77.0	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	877-09-8	Tetrachloro-m-xylene			29.7	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/21/2009 08	09/29/2009	877-09-8	Tetrachloro-m-xylene [2C]		1	27.1	ug/kg
	CGS-US1-091709	RSI0634-01	09/17/2009	09/18/2009 09	09/18/200	STL00234	Percent Solids		1	27	%
Upstream Duplicate	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC		-	530	ua/ka
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC [2C]			560	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	319-85-7	beta-BHC [2C]			200	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2004	319-85-7	beta-BHC			100	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	319-86-8	delta_BHC			26	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2000	310 96 9	delta BHC [2C]		J	30	ug/kg
	CGS-USD1-091709	PSI0634.02	09/17/2009	09/21/2009 00	09/29/200	519-00-0			1	39	ug/kg
	CCS USD1-091709	RS10034-02	09/17/2009	09/21/2009 08	09/29/200:	56-69-9	gamma-BHC (Lindane)		J	19	ug/kg
	CGS-05D1-091709	RS10634-02	09/17/2009	09/21/2009 08	09/29/2009	58-89-9	gamma-BHC (Lindane) [2C]		J	28	ug/kg
	000 0001-091/09	KS10634-02	09/1//2009	09/21/2009 08	09/29/2009	2051-24-3	Decachlorobiphenyl			37.2	ug/kg
	CGS-USD1-091709	KS10634-02	09/17/2009	09/21/2009 08	09/29/2009	2051-24-3	Decachlorobiphenyl [2C]			45.1	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	877-09-8	Tetrachloro-m-xylene			29.1	ug/kg
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/21/2009 08	09/29/2009	877-09-8	Tetrachloro-m-xylene [2C]			26.7	ug/ka
	CGS-USD1-091709	RSI0634-02	09/17/2009	09/18/2009 09	09/18/2009	STL00234	Percent Solids			27	%
Downstream	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC			210	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2009	319-84-6	alpha-BHC [2C]			240	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2009	319-85-7	beta-BHC [2C]			69	ua/ka
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/200	319-85-7	beta-BHC			73	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2004	319-86-8	delta-BHC		J	32	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2004	319-86-8	delta-BHC [2C]		ĭ	34	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2004	58-89-9	gamma-BHC (Lindane)		·		
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 09	09/20/200	58.80.0	gamma-BHC (Lindana) [201			ND	ug/kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2000 00	09/20/200	2051.24.2	Decachlorobinhand			127.0	ug/Kg
	CGS-DS1-091709	RSI0634-03	09/17/2009	09/21/2009 08	09/29/2009	2051-24-3	Decachiorobiphenyl			27.0	ug/kg
	CGS-DS1-091709	RSI0634-03	00/17/2009	00/21/2009 08	09/29/2009	2001-24-3	Decachiorobiphenyl [2C]			40.9	ug/kg
	CCS DS1-091709	NSI0034-03	09/17/2009	09/21/2009 08	09/29/2009	0//-09-8	retrachioro-m-xylene			28.9	ug/kg
	000-001-091/09	RSI0034-03	09/17/2009	09/21/2009 08	09/29/2009	8/1-09-8	etrachloro-m-xylene [2C]			26.7	ug/kg
	CGS-DS1-091709	KS10634-03	09/17/2009	09/18/2009 09	09/18/200	STL00234	Percent Solids			27	%
Downstroom Du V		DOI000 1 D	00/17								
Downstream Duplicate	CGS-DSD1-091709	KSI0634-04	09/17/2009	09/21/2009 08	09/29/200	319-84-6	alpha-BHC			93	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/200	319-84-6	alpha-BHC [2C]			110	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/200	319-85-7	beta-BHC [2C]			180	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/2009	319-85-7	beta-BHC			170	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/200	319-86-8	delta-BHC		1	24	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/2004	319-86-8	delta-BHC [2C]		ř ł	26	
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 09	09/29/2004	58,89 0	damma_BHC (Lindone)		-	ND	ug/kg
	CGS-DSD1-091709	RSI0634.04	09/17/2009	09/21/2009 08	00/20/200	50.00-0	gamma BHO (Lindane)				ug/kg
	CGS-DSD1-091709	PSI0624 04	00/17/2009	00/21/2009 08	00/20/200	00-09-9	gamma-Bric (Lindane) [2C]			ND	ug/kg
	CCS DSD1-091709	RS10034-04	09/17/2009	09/21/2009 08	09/29/2009	2051-24-3	Decachiorobiphenyl			21.2	ug/kg
	000 0001-091/09	RSI0034-04	09/17/2009	09/21/2009 08	09/29/200	2051-24-3	Decachlorobiphenyl [2C]			34.9	ug/kg
	000 0001-091/09	KS10634-04	09/17/2009	09/21/2009 08	09/29/200	8/7-09-8	Tetrachloro-m-xylene			20.7	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/21/2009 08	09/29/200	877-09-8	Tetrachloro-m-xylene [2C]			19.1	ug/kg
	CGS-DSD1-091709	RSI0634-04	09/17/2009	09/18/2009 09	09/18/2009	STL00234	Percent Solids			38	%

SDG: RSI0634 – Test America Laboratories (STL), Amherst, NY

Deliverables

The data packages as submitted to Olin Corporation are complete as stipulated under the Quality Assurance Project Plan (QAPP) for United States Environmental Protection Agency (USEPA) Methods 8081A and percent solids by ASTM D2216-90.

Sample Integrity

Samples within this sample delivery group (SDG) were submitted to the Test America laboratory in Amherst, NY (Buffalo) for chlorinated pesticide analyses. The sample cooler temperature measured 2.0°C at the laboratory which is within the required limit of 4°C \pm 2°. The proper bottles and preservatives were used, the Chain of Custody was properly relinquished, and the correct analytical methods were employed.

Sample Identification

This SDG contains the following water, soil and quality control (QC) samples collected on September 17, 2009:

SDG RSI0634

Sample ID	Sample ID	Sample ID	Sample ID
US-1-091709	USD-1-091709	DS-1-091709	DSD-1-091709

Chlorinated Pesticides (8081A)

The samples in this SDG were submitted for chlorinated pesticides by USEPA Method 8081A.

Holding Times

The extraction and analytical logs indicate that applicable holding times were met for samples submitted for chlorinated pesticide analyses.

Practical Quantitation Limits

The practical quantitation limits (PQLs) as stipulated in the QAPP were not met for all sediment samples submitted for the analysis of chlorinated pesticides by USEPA Method 8081A. Due to matrix interferences samples US-1-091709, USD-1-091709, DS-1-091709 and DSD-1-091709 required a 10X dilution.

Calibration

The initial and continuing calibration data for this SDG indicates that the applicable calibration criteria were met for samples submitted for chlorinated pesticide analyses. The column breakdown for Endrin and DDT was assessed each day the Olin samples were analyzed and the percent degradation of Endrin and DDT were within QC limits.

Surrogates

The surrogate recoveries were within applicable QC limits as stipulated by the laboratory for samples submitted for pesticide analysis.

Internal Standards

The internal standard (IS) recoveries were within applicable QC limits as stipulated by the laboratory for volatile analysis. No additional qualification of the data was required.

Blank Summary

The analytical results of the laboratory method blanks indicate that chlorinated pesticides were not detected.

Laboratory Control Sample and Standard Reference Material Check

The laboratory control sample (LCS) (ongoing precision and recovery [OPR] sample) spike recoveries and the standard reference material (SRM) check were within the applicable QC advisory limits as specified in the QAPP.

Matrix Spike/Matrix Spike Duplicate

No project samples were selected in the field or by the laboratory for MS/MSD analysis.

Sampling Accuracy

The data was within applicable QC advisory limits; therefore no qualification was required.

Laboratory Duplicate Samples

No samples were selected by the laboratory for duplicate analyses.

Field Duplicate Samples

Samples US-1-091709/USD-1-091709 and DS-1-091709/DSD-1-091709 were selected in the field for duplicate analyses. The relative percent difference (RPD) exceeded QC limits for α -BHC in sample US-1-091709/USD-1-091709. The corresponding results was flagged "J". The RPD for δ -BHC and γ -BHC were within QC limits.

The RPD for α -BHC and β -BHC exceeded QC limits in sample DS-1-091709/DSD-1-091709. The corresponding results were flagged "J". The RPD for γ -BHC could not be assessed since the compound was not detected. The RPD for δ -BHC was within QC limits.

Qualification Table

Data Flag: J = Estimated quantitation based upon QC data

Sample ID	Constituent	Data Flag	Sample ID	Constituent	Data Flag
US-1-091709	α-BHC	J	DS-1-091709	α-BHC	J
				β-BHC	J

Percent Solids

The following criteria were used to evaluate samples submitted for percent solids analysis:

Solids samples with <50% solid content require qualification; if the solid content is <50% but \geq 10%, qualify positive results as estimated "J" and qualify non-detect results as estimated "UJ". The results for

samples US-1-091709/USD-1-091709 and DS-1-091709/DSD-1-091709 contained <50% but ≥10% percent solids and were flagged "J".

Qualification Table

Data Flag: J = Estimated quantitation based upon QC data	а
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Sample ID	Constituent	Data Flag	Sample ID	Constituent	Data Flag
US-1-091709	Percent Solids	J	DS-1-0917091	Percent Solids	J
USD-1-091709	Percent Solids	J	DSD-1-091709	Percent Solids	J

Overall Site Evaluation and Professional Judgment Flagging Changes

The data within this SDG were compared to site data and edits to the DQE flags were required based on professional judgment. Monitoring period completeness, which is the percentage of analytical results judged to be valid, including estimated values, was 100 percent for the September 2009 sampling event. Typically, project objectives are met when completeness is 90 percent or better.

James E. Young Date: October 26, 2009 Prepared by:

<u>TestAmerica</u>

THE LEADER IN ENVIRONMENTAL TESTING

TestAmerica Laboratories, Inc.

ANALYTICAL REPORT

Work Order #: RSI0634

Project#: NY3A9025 Site Name: <u>OLIN CORPORATION</u> Task: Charles Gibson Site

Mr. Mike Bellotti Olin Corporation Environmental Remediation Group 3855 North Ocoee Street, Suite 200 Cleveland, TN 37312

CC: Mr. Michael Walker

TestAmerica Laboratories

Brian J. Fischer Project Manager

Muy

Mary Ann Neary Analyst

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THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Received:

09/17/09 Reported: 10/02/09 13:42

2/275

Project: Olin - Charles Gibson site OLN Project Number:

TestAmerica Buffalo Current Certifications

As of 1/27/2009

STATE	Program	Cert # / Lab ID
Arkansas	CWA, RCRA, SOIL	88-0686
California*	NELAP CWA, RCRA	01169CA
Connecticut	SDWA, CWA, RCRA, SOIL	PH-0568
Florida*	NELAP CWA, RCRA	E87672
Georgia*	SDWA, NELAP CWA, RCRA	956
Illinois*	NELAP SDWA, CWA, RCRA	200003
lowa	SW/CS	374
Kansas*	NELAP SDWA, CWA, RCRA	E-10187
Kentucky	SDWA	90029
Kentucky UST	UST	30
Louisiana*	NELAP CWA, RCRA	2031
Maine	SDWA, CWA	N Y0044
Maryland	SDWA	294
Massachusetts	SDWA, CWA	M-NY044
Michigan	SDWA	9937
Minnesota	SDWA,CWA, RCRA	036-999-337
New Hampshire*	NELAP SDWA, CWA	233701
New Jersey*	NELAP, SDWA, CWA, RCRA,	NY455
New York*	NELAP, AIR, SDWA, CWA, RCRA, CLP	10026
Oklahoma	CWA, RCRA	9421
Pennsylvania *	NELAP CWA,RCRA	68-00281
Tennessee	SDWA	02970
Texas *	NELAP CWA, RCRA	T104704412-08-TX
USDA	FOREIGN SOIL PERMIT	S-41579
USDOE	Department of Energy	DOECAP-STB
Virginia	SDWA	278
Washington*	NELAP CWA,RCRA	C1677
Wisconsin	CWA, RCRA	998310390
West Virginia	CWA,RCRA	252

*As required under the indicated accreditation, the test results in this report meet all NELAP requirements for parameters for which accre ditation is required or available. Any exceptions to NELAP requirements are noted in this report.

TestAmerica Buffalo

10 Hazelwood Drive Amherst, NY 14228 tel 716-691-2600 fax 716-691-7991 www.testamericainc.com



OER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Project: Olin - Charles Gibson site Project Number: OLN

Case Narrative

According to 40CFR Part 136.3, pH, Chlorine Residual, Dissolved Oxygen, Sulfite, and Temperature analyses are to be performed immediately after aqueous sample collection. When these parameters are not indicated as field (e.g. field-pH), they were not analyzed immediately, but as soon as possible after laboratory receipt.

A pertinent document is appended to this report, 1 page, is included and is an integral part of this report.

Reproduction of this analytical report is permitted only in its entirety. This report shall not be reproduced except in full without the written approval of the laboratory.

TestAmerica Laboratories, Inc. certifies that the analytical results contained herein apply only to the samples tested as received by our Laboratory.

Received: Reported:

09/17/09 10/02/09 13:42 THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RS10634

Project: Olin - Charles Gibson site Project Number: OLN

Received: Reported:

09/17/09 10/02/09 13:42

	DATA QUALIFIERS AND DEFINITIONS
D08	Dilution required due to high concentration of target analyte(s)
J QFL	Analyte detected at a level less than the Reporting Limit (RL) and greater than or equal to the Method Detection Limit (MDL). Concentrations within this range are estimated. Florisil clean-up (EPA 3620) performed on extract.
NR	Any inclusion of NR indicates that the project specific requirements do not require reporting estimated values below the laboratory reporting limit.

ADDITIONAL COMMENTS

Results are reported on a wet weight basis unless otherwise noted.

TestAmerica Buffalo 10 Hazelwood Drive Amherst, NY 14228 tel 716-691-2600 fax 716-691-7991 www.testamericainc.com

lestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Project: Olin - Charles Gibson site Project Number: OLN

			Executive	Summa	ry - Detect	ions				
	Sample	Data				Dil	Date	Lab		
Analyte	Result	Qualifiers	RL	MDL	Units	Fac	Analyzed	Tech	Batch	Method
Sample ID: RSI0634-01 (CGS-US1-09	1709 - Solid)			Samp	led: 09	/17/09 11:30	Recvd: 09/17/09 15:00		
Organochlorine Pesticio	des by EPA M	Method 8081A								
alpha-BHC	240	QFL, D08	61	11	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
beta-BHC	260	QFL, D08	61	44	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
delta-BHC	39	QFL, D08,J	61	8.1	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
gamma-BHC (Lindane)	18	QFL, D08,J	61	11	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
General Chemistry Para	meters									
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:15	СЈМ	9118022	Dry Weight
Sample ID: RSI0634-02 (CGS-USD1-0	91709 - Solid)			Samp	led: 09	/17/09 11:40	Rec	/d: 09/17/0	9 15:00
Organochlorine Pesticio	des by EPA M	Method 8081A								
alpha-BHC	530	QFL, D08	61	11	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A
beta-BHC	190	QFL, D08	61	44	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A
delta-BHC	36	QFL, D08,J	61	8.0	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A
gamma-BHC (Lindane)	19	QFL, D08,J	61	11	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A
General Chemistry Para	meters									
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:17	CJM	9118022	Dry Weight
Sample ID: RSI0634-03 (CGS-DS1-09	1709 - Solid)			Samp	led: 09/	17/09 12:00	Recu	/d: 09/17/0	9 15:00
Organochlorine Pesticio	des by EPA M	Method 8081A								
alpha-BHC	210	QFL, D08	60	11	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A
beta-BHC	73	QFL, D08	60	44	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A
delta-BHC	32	QFL, D08,J	60	8.0	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A
General Chemistry Para	meters									
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:19	CJM	9118022	Dry Weight
Sample ID: RSI0634-04 (CGS-DSD1-0	91709 - Solid)			Samp	led: 09/	17/09 12:20	Recv	rd: 09/17/0	9 15:00
Organochlorine Pesticio	des by EPA M	Method 8081A								에는 제품을 가지. 1942년 - 1947년 월
alpha-BHC	93	QFL, D08	43	7.8	ug/kg dry	10.0	09/29/09 04:20	MAN	9120007	8081A
beta-BHC	170	QFL, D08	43	31	ug/kg dry	10.0	09/29/09 04:20	MAN	9120007	8081A
delta-BHC	24	QFL, D08,J	43	5.7	ug/kg dry	10.0	09/29/09 04:20	MAN	9120007	8081A
General Chemistry Para	imeters									
Percent Solids	38		0.010	NR	%	1.00	09/18/09 15:21	CJM	9118022	Dry Weight
					1					

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Received: 09/17/09 Reported: 10/02/09 13:42



THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Clevela	and, TN	Work Order: RSI0634			Received:	09/17/09
3855 North Ocoee St. Suite 200					Reported:	10/02/09 13:42
Cleveland, TN 37312		Project: Olin - Charles	Gibson site			
		Project Number: C	DLN			

Sample Summary

Sample Identification	Lab Number	Client Matrix	Date/Time Sampled	Date/Time Sample Received Qualifiers
CGS-US1-091709	RSI0634-01	Solid	09/17/09 11:30	09/17/09 15:00
CGS-USD1-091709	RS10634-02	Solid	09/17/09 11:40	09/17/09 15:00
CGS-DS1-091709	RS10634-03	Solid	09/17/09 12:00	09/17/09 15:00
CGS-DSD1-091709	RS10634-04	Solid	09/17/09 12:20	09/17/09 15:00



THE LEADER IN ENVIRONMENTAL TESTING

Received: 09/17/09 Reported: 10/02/09 13:42

Olin Chlor Alkali Products - Cleveland, TN 3855 North Occee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Project: Olin - Charles Gibson site Project Number: OLN

			4	Analytica	Report					
Analyte	Sample Result	Data Qualifiers	RL	MDL	Units	Dil Fac	Date Analyzed	Lab Tech	Batch	Method
Sample ID: RSI0634-01 (CGS-US1-09	1709 - Solid)			Samp	oled: 09/	17/09 11:30	Recu	/d: 09/17/0	9 15:00
Organochlorine Pesticio	des by EPA M	lethod 8081/	\							
alpha-BHC	240	QFL, D08	61	11	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
beta-BHC	260	QFL, D08	61	44	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
delta-BHC	39	QFL, D08, J	61	8.1	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
gamma-BHC (Lindane)	18	QFL, D08,J	61	ji (. 11)	ug/kg dry	10.0	09/29/09 02:32	MAN	9120007	8081A
Decachlorobiphenyl	169 %	QFL, D08	Surr Limits:	(42-146%)			09/29/09 02:32	MAN	9/20007	8081A
Tetrachloro-m-xylene	121 %	QFL, D08	Surr Limits:	(37-135%)			09/29/09 02:32	MAN	9120007	8081A
General Chemistry Para	meters									
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:15	СЈМ	9118022	Dry Weight

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THE LEADER IN ENVIRONMENTAL TESTING

09/17/09

10/02/09 13:42

Received:

Reported:

Olin Chlor Alkali Products - Cleveland, TNWork Order: RSI06343855 North Ocoee St. Suite 200Cleveland, TN 37312Project: Olin - Charles Gibson site

Project Number: OLN

		Analytical Report									
Analyte	Sample Result	Data Qualifiers	RL	MDL	Units	Dil Fac	Date Analyzed	Lab Tech	Batch	Method	
Sample ID: RSI0634-02 (CG	S-USD1-0	91709 - Solid) •		Samp	led: 09/	17/09 11:40	Rec	vd: 09/17/0	9 15:00	
Organochlorine Pesticides	by EPA N	Method 8081/	\								
alpha-BHC	530	QFL, D08	61	11	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A	
beta-BHC	190	QFL, D08	61	44	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A	
delta-BHC	36	QFL, D08,J	61	8.0	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A	
gamma-BHC (Lindane)	19	QFL, D08,J	61	11	ug/kg dry	10.0	09/29/09 03:08	MAN	9120007	8081A	
Decachlorobiphenyl	154 %	QFL, D08	Surr Limits:	(42-146%)			09/29/09 03:08	MAN	9/20007	8081A	
Tetrachloro-m-xylene	120 %	QFL, D08	Surr Limits:	(37-135%)			09/29/09 03:08	MAN	9/20007	8081A	
General Chemistry Parame	ters										
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:17	CJM	9118022	Dry Weight	

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THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN

9/275

Received: 09/17/09 Reported: 10/02/09 13:42

3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Project: Olin - Charles Gibson site Project Number: OLN

			4	Analytical	Report						
Analyte	Sample Result	Data Qualifiers	RL	MDL	Units	Dil Fac	Date Analyzed	Lab Tech	Batch	Method	
Sample ID: RSI0634-03	(CGS-DS1-09	1709 - Solid)			Samp	led: 09	17/09 12:00	Recy	/d: 09/17/(9 15:00	
Organochlorine Pestic	ides by EPA I	Method 8081/	7						estations) States estation		
alpha-BHC	210	QFL, D08	60	11	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A	
beta-BHC	73	QFL, D08	60	44	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A	
delta-BHC	32	QFL, D08,J	60	8.0	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A	
gamma-BHC (Lindane)	ND	QFL, D08	60	10	ug/kg dry	10.0	09/29/09 03:44	MAN	9120007	8081A	
Decachlorobiphenyl	112 %	QFL, D08	Surr Limits:	(42-146%)			09/29/09 03:44	MAN	9/20007	8081A	
Tetrachloro-m-xylene	120 %	QFL, D08	Surr Limits:	(37-135%)			09/29/09 03:44	MAN	9120007	8081A	
General Chemistry Par	ameters					an ting. An Arts					
Percent Solids	27		0.010	NR	%	1.00	09/18/09 15:19	СЈМ	9118022	Dry Weight	

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THE LEADER IN ENVIRONMENTAL TESTING

 Olin Chlor Alkali Products - Cleveland, TN
 Work Order: RSI0634
 Received: 09/17/09

 3855 North Ocoee St. Suite 200
 Reported: 10/02/09 13:42

 Cleveland, TN 37312
 Project: Olin - Charles Gibson site

 Project Number:
 OLN

				Analytical I	Report		
Analyte	Sample Result	Data Qualifiers	RL	MDL	CUnitsF	il Date ac Analyzed	Lab Tech Batch Method
Sample ID: RSI0634-04 (CC	S-DSD1-	091709 - Solia	J)		Sampled	09/17/09 12:20	Recvd: 09/17/09 15:00
Organochlorine Pesticide	s by EPA	Method 8081	<u>A</u>				
alpha-BHC beta-BHC delta-BHC gamma-BHC (Lindane)	93 170 24 ND	QFL, D08 QFL, D08 QFL, D08,J QFL, D08	- 43 43 43 43	7.8 31 5.7 7.5	ug/kg dry 10 ug/kg dry 10 ug/kg dry 10 ug/kg dry 11	0.0 09/29/09 04:20 0.0 09/29/09 04:20 0.0 09/29/09 04:20 0.0 09/29/09 04:20 0.0 09/29/09 04:20 0.0 09/29/09 04:20) MAN 9120007 8081A) MAN 9120007 8081A) MAN 9120007 8081A) MAN 9120007 8081A
Decachlorobiphenyl Tetrachloro-m-xylene	122 % 119 %	QFL, D08 QFL, D08	Surr Limits: Surr Limits:	(42-146%) (37-135%)		09/29/09 04:20 09/29/09 04:20) MAN 9120007 8081A) MAN 9120007 8081A
General Chemistry Parame Percent Solids	eters 38		0.010	NR	% 1	00 09/18/09 15:21	I CJM 9I18022 Dry Weight



THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312

Work Order: RSI0634

Project: Olin - Charles Gibson site Project Number: OLN

SAMPLE EXTRACTION DATA

			Wt/Vol		Extract			Lab	하는 것은 것이 가 있었다. 이야지 같은 것은 것은 것은 것은 것이 있는 것이다.
Parameter	Batch	Lab Number	Extracte	Units	Volume	Units	Date Prepared	Tech	Extraction Method
General Chemistry Parameters									
Dry Weight	9118022	RSI0634-01	10.00	g	10.00	g	09/18/09 09:20	CJM	Dry Weight
Dry Weight	9118022	RSI0634-02	10.00	g	10.00	g	09/18/09 09:20	СЈМ	Dry Weight
Dry Weight	9118022	RSI0634-03	10.00	g	10.00	g	09/18/09 09:20	CJM	Dry Weight
Dry Weight	9118022	RS10634-04	10.00	g	10.00	g	09/18/09 09:20	CJM	Dry Weight
Organochlorine Pesticides by EP	A Method 8	081A							이 지수 같은 것이다.
8081A	9120007	RSI0634-02	30.15	g	10.00	mL	09/21/09 08:00	СХМ	3550B GC
8081A	9120007	RSI0634-01	30.28	g	10.00	mL	09/21/09 08:00	CXM	3550B GC
8081A	9120007	RSI0634-04	30.44	g	10.00	mL	09/21/09 08:00	CXM	3550B GC
8081A	9120007	RSI0634-03	30.54	g	10.00	mL	09/21/09 08:00	СХМ	3550B GC
8081A	9120007	RSI0634-03	30.54	g	10.00	mL	09/21/09 08:00	СХМ	3550B GC

Received: 09/17/09 Reported: 10/02/09 13:42 TestAmerica

THE LEADER IN ENVIRONMENTAL TESTING

Olin Chlor Alkali Products - Cleveland, TN 3855 North Ocoee St. Suite 200 Cleveland, TN 37312 Work Order: RSI0634

Received: 09/17/09

Reported: 10/02/09 13:42

Project: Olin - Charles Gibson site Project Number: OLN

			L	BORATOR	QC DATA					
	Source	Spike					%	% REC	% RPD	Data
Analyte	Result	Level	RL	MDL	Units	Result	REC	Limits	RPD Limit	Qualifiers
Organochlorine Pesticide	s by EPA	Method 80	<u>)81A</u>							
Blank Analyzed: 09/25/09	(Lab Num	nber:91200	07-BLK1, E	latch: 9120007)						
alpha-BHC			1.7	0.30	ug/kg wet	ND				
alpha-BHC [2C]			1.7	0.30	ug/kg wet	ND				
beta-BHC			1.7	1.2	ug/kg wet	ND				
beta-BHC [2C]			1.7	1.2	ug/kg wet	ND				
delta-BHC			1.7	0.22	ug/kg wet	ND				
delta-BHC [2C]			1.7	0.22	ug/kg wet	ND				
gamma-BHC (Lindane)			1.7	0.29	ug/kg wet	ND				
gamma-BHC (Lindane)			1.7	0.29	ug/kg wet	ND				
[2C]				la de la composición	a shekara sheka					
Surrogate:					ug/kg wet		92	42-146		
Decachlorobiphenyl Surrogate: Decachlorobinhenyl (2C)					ug/kg wet		111	42-146		
Surrogate: Tetrachloro-m-xvlene					ug/kg wet		79	37-135		
Surrogate: Tetrachloro-m-xylene					ug/kg wet		54	37-136		
LCS Analvzed: 09/25/09 (Lab Numb	er:912000	7-BS1. Bate	:h: 9 20007)			an a			
alpha-BHC		16	1.6	0.29	ua/ka wet	14.9	91	49-120		
alpha-BHC [2C]		16	1.6	0.29	ug/kg wet	13.1	80	49-120		
beta-BHC		16	1.6	1.2	ug/kg wet	16.7	102	56-120	n se se ga se Se segne se fina	
beta-BHC [2C]		16	1.6	1.2	ug/kg wet	14.6	89	56-120		
delta-BHC		16	1.6	0.22	ug/kg wet	16.0	98	45-123		
delta-BHC [2C]		16	1.6	0.22	ug/kg wet	14.8	90	45-123		
gamma-BHC (Lindane)		16	1.6	0.28	ug/kg wet	15.6	96	50-120		
gamma-BHC (Lindane) [2C]		16	1.6	0.28	ug/kg wet	13.8	84	50-120		
Surrogate:					ug/kg wet		97	42-146		
Surrogate: Decachlorobiphenvl [2C]					ug/kg wet		110	42-146		
Surrogate: Tetrachloro-m-xvlene					ug/kg wet		108	37-135		
Surrogate: Tetrachloro-m-xylene					ug/kg wet		73	37-136		

TestAmerica Buffalo

10 Hazelwood Drive Amherst, NY 14228 tel 716-691-2600 fax 716-691-7991 www.testamericainc.com

Chain of		Temperatu	re on Peceip		ř	∋stA	merico	
Custody Record		Drinking M	later? Yes C	Epon (1	LEADER IN E	AVIRONMENTAL TESTIN	
Ohn Con .		HUKS War	li otto	3414			Date 9-17-09	Chain of Custody Number
2855 North Acoust ST.	Sult 200	Telephone N	3 L - 15C	OFFAX NUMB	* 5. 5	4166	Lab Number	0 + + + + + + + + + + + + + + + + + + +
ALAND TO THE TO CO	81 L	NAL CONTRES	rich	24	c heo		Analysis (Attach list if Thore space is needed)	
Project Name application (Spice) Unspec Palls	N.L.	Carriav/W/ry/	vit Number					
Contractifications Orden Ouche No.	 -		Matrix	Paris Paris	Itainera di servatives	ç)		opecal instructions Conditions of Placeipt
Sample I.D. No. and Description (Containers for each sample may be combined on one line)	Date	Time }	Pape Dabe Moontry	HINCE HISCON FURMER	WPOH TOVE	99		
CGS. USI . 09 17 69	6-17-0	1130	K			X		HAN T KI
CGS . USDI . 021 . 091		ohii	7			×		
C65 . 151 . 091709		10.01	*	-		12		
CG3. NSDI. CALTON	≯	1 acro	. 7			X		
Possible Hazard Identification	Poison B	Nitmourn S	Imple Disposal Return To Chien	D Duepo	1 444 A	Achine For	Adonths toporer then 1 a	essessed if samples are relatined working
Tum Around Time Acquired	21 Days	1 Sumo	Comment.	- J OC Field	unimuraturits (Spe	<u>؟</u>		
Unter the wall		0-1-1	13100	1. Abcs	and aller	TH J	Buttan	Daviality 1 1 2m
2. Rethrausted By		Date	entit	2 Aeca	言を			
3. Retirquisted By			1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	J. Pase	(read By			Date Time
Comments						۱۲ 		
OUSTRUBUTION: WHITE - Returned to Chevy with Report, CA	NARV - Util	th the Sample.	VINK · Field Copy			*		
						7	Ś	

TestAmerica Buffalo SDG: CLASS: PEST METHOD: 8081A

ANALYSES DATA PACKAGE COVER PAGE

8081A

Laboratory: TestAmerica Buffalo

Client: Olin Chlor Alkali Products - Cleveland, TN

Client Sample Id: CGS-US1-091709 CGS-USD1-091709 CGS-DS1-091709 CGS-DSD1-091709

SDG:

Project: Olin - Charles Gibson site - NY3A9025AE03759

Lab Sample Id: <u>RSI0634-01</u> <u>RSI0634-02</u> <u>RSI0634-03</u>

RSI0634-04

Form 2 SURROGATE STANDARD RECOVERY AND RT SUMMARY

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Sequence:	<u>RI92924</u>	Instrument:	<u>HP6890-6</u>
Matrix:	Solid	Calibration:	<u>R9H1803</u>

Surrogate Compound	Spike Level ug/kg wet	% Recovery	Recovery Limits	RT	Calibration Mean RT	RT Diff	RT Diff Limit	Q
Blank (9120007-BLK1)			Lab File ID: 6/	A46193	Analyzed	: 09/25/09 12:0	9	
Decachlorobiphenyl	6.62	92	42 - 146	22.31588	22.68772	-0.3718	+/-1.0	
Decachlorobiphenyl [2C]	6.62	111	42 - 146	25.5428	26.11406	-0.5713	+/-1.0	
Tetrachloro-m-xylene	6.62	79	37 - 135	9.358156	9.636484	-0.2783	+/-1.0	
Tetrachloro-m-xylene [2C]	6.62	54	37 - 136	10.71218	11.02601	-0.3138	+/-1.0	
LCS (9I20007-BS1)			Lab File ID: 6/	A46181	Analyzed: 09/25/09 04:58			
Decachlorobiphenyl	6.54	97	42 - 146	22.31062	22.68772	-0.3771	+/-1.0	
Decachlorobiphenyl [2C]	6.54	110	42 - 146	25.53448	26.11406	-0.5796	+/-1.0	
Tetrachloro-m-xylene	6.54	108	37 - 135	9.357999	9.636484	-0.2785	+/-1.0	
Tetrachloro-m-xylene [2C]	6.54	73	37 - 136	10.71147	11.02601	-0.3145	+/-1.0	

Form 2

SURROGATE STANDARD RECOVERY AND RT SUMMARY

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Sequence:	<u>R192925</u>	Instrument:	<u>HP6890-6</u>
Matrix:	Solid	Calibration:	<u>R9I2904</u>

Surrogate	Spike	% Baaawaa	Recovery	рт	Calibration	PT D:ff	RT Diff	
Compound	Level ug/kg dry	Recovery		KI	Mean K1	KI DIII		
CGS-US1-091709 (RSI0634-01)			Lab File ID: 6.	A46240	Analyzed	: 09/29/09 02:3	2	
Decachlorobiphenyl	24.5	169	42 - 146	22.10082	22.08204	0.0188	+/-1.0	*
Decachlorobiphenyl [2C]	24.5	314	42 - 146	25.25846	25.21592	0.0425	+/-1.0	*
Tetrachloro-m-xylene	24.5	121	37 - 135	9.194136	9.178268	0.0159	+/-1.0	
Tetrachloro-m-xylene [2C]	24.5	110	37 - 136	10.54256	10.53044	0.0121	+/-1.0	
CGS-USD1-091709 (RSI0634-02)			Lab File ID: 6	A46241	Analyzed	: 09/29/09 03:0	8	
Decachlorobiphenyl	24.2	154	42 - 146	22.0966	22.08204	0.0146	+/-1.0	*
Decachlorobiphenyl [2C]	24.2	186	42 - 146	25.24903	25.21592	0.0331	+/-1.0	*
Tetrachloro-m-xylene	24.2	120	37 - 135	9.192714	9.178268	0.0144	+/-1.0	
Tetrachloro-m-xylene [2C]	24.2	110	37 - 136	10.54188	10.53044	0.0114	+/-1.0	
CGS-DS1-091709 (RSI0634-03)			Lab File ID: 6	A46242	Analyzed	: 09/29/09 03:4	4	
Decachlorobiphenyl	24.1	112	42 - 146	22.095	22.08204	0.0130	+/-1.0	
Decachlorobiphenyl [2C]	24.1	169	42 - 146	25.24854	25.21592	0.0326	+/-1.0	*
Tetrachloro-m-xylene	24.1	120	37 - 135	9.187011	9.178268	0.0087	+/-1.0	
Tetrachloro-m-xylene [2C]	24.1	111	37 - 136	10.53767	10.53044	0.0072	+/-1.0	
CGS-DSD1-091709 (RSI0634-04)			Lab File ID: 6.	A46243	Analyzed	: 09/29/09 04:2	0	
Decachlorobiphenyl	17.3	122	42 - 146	22.0946	22.08204	0.0126	+/-1.0	
Decachlorobiphenyl [2C]	17.3	202	42 - 146	25.24826	25.21592	0.0323	+/-1.0	*
Tetrachloro-m-xylene	17.3	119	37 - 135	9.188422	9.178268	0.0102	+/-1.0	
Tetrachloro-m-xylene [2C]	17.3	110	37 - 136	10.53632	10.53044	0.0059	+/-1.0	

Form 3

LCS / LCS DUPLICATE RECOVERY

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	<u>Olin - Charles Gibson site - NY3A9025AE03759</u>
Matrix:	Solid	Spike standard:	<u>9080196</u>
Batch:	<u>9120007</u>	Laboratory ID:	<u>9I20007-BS1</u>
Preparation:	<u>3550B GC</u>	Initial/Final:	<u>30.56 g / 10 mL</u>

COMPOUND	SPIKE ADDED	UNITS	LCS CONCENTRATION	LCS % REC. #	QCLIMITS REC.
alpha-BHC	16.4	ug/kg wet	14.9	91	49 - 120
alpha-BHC [2C]	16.4	ug/kg wet	13.1	80	49 - 120
beta-BHC	16.4	ug/kg wet	16.7	102	56 - 120
beta-BHC [2C]	16.4	ug/kg wet	14.6	89	56 - 120
delta-BHC	16.4	ug/kg wet	16.0	98	45 - 123
delta-BHC [2C]	16.4	ug/kg wet	14.8	90	45 - 123
gamma-BHC (Lindane)	16.4	ug/kg wet	15.6	96	50 - 120
gamma-BHC (Lindane) [2C]	16.4	ug/kg wet	13.8	84	50 - 120

Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

PREPARATION BATCH SUMMARY

8081A

Laborat	ory: <u>TestAmerica E</u>	Buffalo	SDG:				
Client:	Olin Chlor Alk	ali Products - Cleveland, TN	Project	<u>Olin</u>	Olin - Charles Gibson site - NY3A9025AE03759		
Batch:	9120007	Batch Matrix: Solid	Prepara	tion: <u>3550</u>	<u>B GC</u>		
	SAMPLE NAME	LAB SAMPLE ID	INITIAL	FINAL	DATE PREPARED	TOT/DIS	
	Blank	9I20007-BLK1	30.2 g	10.0 mL	09/21/09 08:00	N/A	
	LCS	9I20007-BS1	30.6 g	10.0 mL	09/21/09 08:00	N/A	
	CGS-US1-091709	RSI0634-01	30.3 g	10.0 mL	09/21/09 08:00	N/A	
	CGS-USD1-091709	RSI0634-02	30.2 g	10.0 mL	09/21/09 08:00	N/A	
	CGS-DS1-091709	RSI0634-03	30.5 g	10.0 mL	09/21/09 08:00	N/A	
	CGS-DSD1-091709	RSI0634-04	30.4 g	10.0 mL	09/21/09 08:00	N/A	

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METHOD DETECTION AND REPORTING LIMITS

8081A

Laboratory:	TestAmerica Buffalo
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Client: Olin Chlor Alkali Products - Cleveland, TN

Matrix: Solid

SDG:

Project: Olin - Charles Gibson site - NY3A9025AE(

<u>olid</u> Instrument: <u>HP68</u>				
Analyte	MDL	MRL	Units	
alpha-BHC	0.30	1.7	ug/kg	
alpha-BHC [2C]	0.30	1.7	ug/kg	
beta-BHC	1.2	1.7	ug/kg	
beta-BHC [2C]	1.2	1.7	ug/kg	
delta-BHC	0.22	1.7	ug/kg	
delta-BHC [2C]	0.22	1.7	ug/kg	
gamma-BHC (Lindane)	0.29	1.7	ug/kg	
gamma-BHC (Lindane) [2C]	0.29	1.7	ug/kg	

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Form 1 ORGANIC ANALYSIS DATA SHEET

CGS-US1-091709

8081A

Laboratory:	TestAmerica Buffalo				SDG:			
Client:	Olin Chlor Alkali	i Products - Clevelan	<u>d, TN</u>		Project:	<u>Olin - Charles Gi</u>	bson site - NY3A90	25AE03759
Matrix:	<u>Solid</u>	Laborato	ry ID:	<u>RSI0634</u>	-01	File ID:	<u>6A46240</u>	
Sampled:	<u>09/17/09 11:30</u>	Prepared	:	<u>09/21/09</u>	08:00	Analyzed:	09/29/09 02:32	
Solids:	26.93	Preparati	on:	3550B C	C	Initial/Final:	<u>30.28 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>R192925</u>		Calibration:	<u>R912904</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CON	C. (ug/kg)	Q
319-84-6	alpha-BHC				10	240		D
319-85-7	beta-BHC				10	260		D
319-86-8	delta-BHC				10	39		л
58-89-9	gamma-BHC (Lii	ndane)			10		18	JPD
SYSTEM MONITORING COMPOUND		ADDED ((ug/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q	
Decachlorobiphenyl 24			24.5	5	41.3	169 42 - 146		*
Tetrachloro-m-xy	lene		24.5	5	29.7	121	37 - 135	l

* Values outside of QC limits

Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:51:11
Operator	: tchrom	Sample Name	: RSI0634-01
Sample Number	:	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/40
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 02:32:22	Cycle	: 14

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46240.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46240.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46240.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46240.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46240.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.43	7870		в	0.00787	785.89
2	7.67	6958		В	0.00696	1758.60
3	7.78	21337		V	0.02134	6852.48
4	8.82	2352		В	0.00235	980.47
5	9.10	16956		В	0.01696	5353.37
6	9.19	39045	Tetrachloro-m-xylene	V	0.00242	12363.80
7	9.37	2308	•	В	0.00231	835.05
8	9.96	8856		В	0.00886	1960.24
9	10.29	21079		В	0.02108	6751.52
10	10.76	729758	alpha-BHC	В	0.01956	239800.79
11	11.20	5585		В	0.00558	730.16
12	11.61	12127	gamma-BHC	В	0.00150	3680.38
13	11.86	311297	beta-BHC	В	0.02139	94649.54
14	12.30	46039	delta-BHC	В	0.00318	14404.71
15	12.47	38638	<u> </u>	В	0.03864	12080.67
16	12.80	15714	Heptachlor-NUF	В	7.04e-04	4570.86

09/29/2009 07:51:11 Result: H:\TURBO6\6890-06\6-SEQ46\6A46240.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	13 30	2050		B	0.00206	1150 35
18	13.52	5094	Aldrin N	IFB	0.00200	1046 33
19	13 00	22253		, C, D	0.00107	4991 88
20	14 18	15644		B	0.02220	4096 21
20	1/ 31	0000			0.01004	2155 34
27	14.63	160531			0.00331	12153.04
22	15.01	7256	Hent enovide		1.010-04	000 27
23	15.01	10028	Tiehr ehovide		0 01003	2855.82
24	15.28	32713	aamma chlordano		0.01093	2000.02
20	15.20	04525	yanina choruane		0.00100	22566 21
20	15.79	25463			0.09432	6010.21
21	15 70	25405			0.02040	5620.27
20	15.73	10901	Fadoculton I	łů	5 410 04	2472 22
29	16.02	12021	Endosullan		0.410-04	24/3.22
21	16.03	12199			0.01200	3140.73
20	10.21	40071	Dialdrin		0.04007	4727.00
22	10.43	4/00 5114	Dielann		0.00109	1131.02
24	10.04	2462			0.00011	1047.20
24	10.73	3402	Endrin		0.00340	1170.00
30	17.04	33040	Enonn		0.00200	11202.09
30	17.04	7 148			0.00715	2270.51
3/	17.10	7910			0.00138	2480.54
30	11.11	44708	4,4-001		0.00442	6202.90
39	10.10	11300	Enduin aldalauda	B	0.01130	2608.14
40	18.20	22964	Endrin aldenyde	l S	0.00158	6509.33
41	10.00	299804			0.29980	76293.88
42	18.79	8805	metnoxycnior		0.00194	2579.84
43	18.69	12000			0.01289	3498.59
44	19.00	234/6/		l N	0.23477	51585.80
40	19.39	95842			0.09584	18248.63
40	19.59	2/0/0			0.02/5/	5822.31
47	19.70	24011	The states to stars a		0.02401	5621.62
40	19.03	31573	Enann ketone		0.00329	6060.12
49	20.05	10050	、		0.00004	1902.04
50	20.17	120011	λ	. / š	0.12001	28039.30
51	20.01	21/19	l l	V 5	0.02178	4843.98
52	20.74	122004		VĽ	0.00107	104004.00
55	21.09	133094			0.13309	30304.02
54	21.00	2433			0.00243	492.23
50	21.01	4933			0.00493	1497.00
50	21.71	3040			0.00304	1000.00
57	22.00	0.0009	Descehlershiphen	ы V	0.00009	10300.03
50	22.10	94799 70755	Decachiorobiphen	yi v	0.00337	10/2/.09
59	22.30	12100 59166			0.07270	13994.90
61	22.30	00100			0.00017	10200.00
62	22.04	21075			0.00033	5960.09
62	23.11	7/070		D	0.03107	13262 25
60	23.10	24019		ם ק	0.07400	10202.00
65	20.99	6705		0 9	0.02420 0.0670	4024.47
05	24.00			D	0.00070	1239.92
		4118493			2.70611	1.07e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98825	Date	: 09/29/2009 07:51:15
Operator	: tchrom	Sample Name	: RSI0634-01
Sample Number	:	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/40
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.00000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 02:32:22	Cycle	: 14

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46240.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46240.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46240.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46240.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46240.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.30	10329		В	0.01033	3762.00
2	9.33	1642		В	0.00164	607.43
5	10.54	15140	Tetrachloro-m-xylene (В	0.00221	4873.81
6	10.92	9950		В	0.00995	2546.24
9	11.91	9596		В	0.00960	3055.66
10	12.16	3239		В	0.00324	618.77
11	12.30	302630	alpha-BHC (2C)	V	0.02166	95050.40
12	13.18	8553	,	В	0.00855	2289.26
13	13.26	10233	gamma-BHC (2C)	V	0.00213	2108.53
14	13.48	136116	beta-BHC (2C)	В	0.02280	39038.00
15	14.23	19514	delta-BHC (2C)	В	0.00341	5916.85
16	14.49	18871	、 ,	В	0.01887	5380.56
17	14.67	2127		В	0.00213	630.99
18	14.94	3748		В	0.00375	797.73
19	15.35	9383		В	0.00938	2810.52
20	15.86	7561		В	0.00756	1738.14

09/29/2009 07:51:15 Result: H:\TURBO6\6890-06\6-SEQ46\6b46240.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
21	16.06	4184		В	0.00418	1467.68
22	16.27	72752		cВ	0.07275	17998.02
23	16.55	3861	Hept. epoxide (2C) N	ŁВ	3.71e-04	965.76
24	16.77	2954		в	0.00295	975.72
25	17.12	55167	1	В	0.05517	9655.79
26	17.35	14140	alpha chlordane (2C)	V	0.00141	4709.82
27	17.77	26411		В	0.02641	5486.04
28	18.00	3985	Dieldrin (2C)	В	0.00126	1153.39
29	18.43	6972		В	0.00697	1102.62
30	18.52	3768		В	0.00377	1192.41
31	18.63	3103		V	0.00310	768.60
33	18.97	4082	4,4'-DDD (2C)	V	0.00161	1325.13
34	19.22	7509	Endosulfan II (2C)	в	8.30e-04	1575.57
35	19.67	6659	4,4'-DDT (2C)	В	0.00321	1652.55
36	19.91	10161	Endrin aldehyde (2C)	В	6.98e-04	2304.69
37	20.05	7285		V	0.00729	1649.84
39	20.45	133031		В	0.13303	31233.81
40	20.59	5714	Endo. Sulfate (2C)	E	0.00273	1579.06
41	20.86	33389		В	0.03339	7461.92
42	20.97	4045		В	0.00405	1529.90
43	21.16	30132	Methoxychlor (2C)	в	0.00822	7223.67
44	21.47	13176		В	0.01318	2815.15
45	21.64	11115		V	0.01112	2083.18
46	21.77	76580	Endrin ketone (2C)	/ V	0.00977	15609.49
48	22.25	44617		В	0.04462	8994.71
49	22.45	1661	V	В	0.00166	464.84
50	22.69	2464		В	0.00246	589.73
51	23.11	265336		В	0.26534	49363.51
52	23.76	45874		В	0.04587	9419.73
53	24.13	11959		В	0.01196	1552.52
54	24.28	7830		V	0.00783	1254.80
55	24.47	6761		V	0.00676	1312.52
56	24.79	75623		V	0.07562	8246.06
57	25.01	13288		V	0.01329	2071.50
58	25.26	71115	Decachlorobiphenyl (2C	; V	0.00628	9469.17
59	25.49	3293		в	0.00329	664.90
60	25.66	39130		V	0.03913	5667.03
61	27.34	39533		В	0.03953	4973.74
		1737292			1.10831	398789.48

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Form 1 ORGANIC ANALYSIS DATA SHEET

CGS-USD1-091709

8081A

Laboratory:	TestAmerica Buffalo				SDG:			
Client:	Olin Chlor Alkali	Products - Cleve	<u>land, TN</u>		Project:	ct: Olin - Charles Gibson site - NY3A9025AE03759		
Matrix:	<u>Solid</u>	Labor	atory ID:	<u>RSI063</u>	<u>4-02</u>	File ID:	<u>6A46241</u>	
Sampled:	<u>09/17/09 11:40</u>	Prepa	red:	<u>09/21/0</u>	<u>9 08:00</u>	Analyzed:	<u>09/29/09 03:08</u>	
Solids:	27.37	Prepa	ration:	<u>3550B (</u>	<u>3C</u>	Initial/Final:	<u>30.15 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>RI92925</u>		Calibration:	<u>R912904</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CON	C. (ug/kg)	Q
319-84-6	alpha-BHC				10		530	D
319-85-7	beta-BHC				10	190		D
319-86-8	delta-BHC			_	10	36		JD
58-89-9	gamma-BHC (Lin	ndane)			10		19	JPD
SYSTEM MONITORING COMPOUND ADDED (ug/) (ug/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q		
Decachlorobiphenyl 24.2			37.2	154	42 - 146	*		
Tetrachloro-m-xy	lene		24	1.2	29.1	120	37 - 135	

* Values outside of QC limits

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98826	Date	: 09/29/2009 07:51:18
Operator	: tchrom	Sample Name	: RSI0634-02
Sample Number	:	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/41
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 03:08:27	Cycle	: 15

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46241.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46241.rst

Inst Method : h:\turbo6\6890-06\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46241.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46241.rst

Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46241.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	7.42	7077		В	0.00708	613.54
2	7.51	1760		В	0.00176	717.31
3	7.67	6918		В	0.00692	1645.85
4	7.77	26019		V	0.02602	8258.40
5	8.81	3310		В	0.00331	1319.76
6	9.10	19338		в	0.01934	6076.96
7	9.19	38394	Tetrachloro-m-xylene	V	0.00240	12037.58
8	9.97	7340	-	В	0.00734	1624.26
9	10.29	16317		В	0.01632	5285.89
10	10.77	1669824	alpha-BHC	В	0.04342	546569.22
11	11.20	4740		В	0.00474	593.04
12	11.61	15281	gamma-BHC	в	0.00158	4707.83
13	11.86	232596	beta-BHC	В	0.01584	70913.24
14	12.30	39264	delta-BHC	В	0.00300	12201.72
15	12.47	35016		В	0.03502	10936.32
16	12.80	15668	Heptachlor	В	7.03e-04	4635.81

09/29/2009 07:51:18 Result: H:\TURBO6\6890-06\6-SEQ46\6A46241.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	13.39	1632		- B	0.00163	931.17
19	13.52	5241	Aldrin MIT-	- B	0.00108	946.22
20	13.99	10264	/	B	0.01026	1986.12
21	14 18	10540	1	B	0.01054	2824.21
22	14 46	42302		B	0.04230	2860 72
23	14 63	143858		v	0 14386	33651 44
24	14 95	1713	Hent enoxide	Ř	1.080-05	647.89
25	15 14	8499		Ř	0.00850	2225 76
26	15.28	27650	namma chlordane	R	0.00000	7320 42
27	15.43	73548	gamma omoraano	v	0.00171	18335.41
28	15 72	13453		Ř	0.01345	3201 29
29	15 79	17346	4 4'-DDE	v	0.0180	4093 38
30	15.91	14484	Endosulfan I	v	6 01e-04	3749 49
31	16.02	9246	Endosunan	v	0.010 04	2411 24
32	16.20	32469		Ř	0.00020	8509.28
33	16.43	3886	Dieldrin	B	0.00247	1426.92
34	16 64	2871	Dicidiiii	B	0.00100	1025 25
35	16.96	36586	Endrin	B	0.00207	12050 20
36	17.04	4941		Ē	0.00207	1715 76
37	17 10	7600	4 4'-DDD	v	0.00404	2203 36
38	17 72	30084		Ř	0.00137	1834 80
30	18 16	50004	ч,ч°001	R	0.00302	1630 17
40	18 25	21784	Endrin aldehyde	v	0.00000	5643.24
40	18 58	230420	chunn aldenyde	Ř	0.00101	58803 81
12	18.80	200420	Methovychlor	B	0.23043	0/0 12
42	18.80	15520	Wothoxychior	v v	0.00141	3670 /1
40	10.00	140073		v	0.01332	28137.00
45	10.00	73445		R	0.14007	13744 27
46	10.50	15158			0.07545	2725.76
40	10.75	15218		1 $\tilde{\lambda}$	0.01510	2670.50
47	10.83	10864	Endrin kotono	$ /\tilde{v} $	0.01022	2921 47
40	20.16	1/0721		/ 6	0.00202	27077 20
50	20.10	18201	Y		0.14072	4226.20
51	20.01	512064			0.01029	121444 20
52	20.75	21833		R	0.01200	5053 <i>1</i> 1
53	21.00	12/501		о В	0.02103	0716 0/
54	21.70	62023			0.12439	11027 61
55	27.55	88003	Decachlorohinhonyl	Ň	0.00292	16206 57
56	22.10	1615666	Decachiotopiphenyi	v	1 61567	129202 42
57	22.40	23025			0.02202	2264 50
58	22.01	23923			0.02393	4525.00
50	23.10	56201		D D	0.02433	4020.99
60	23.00	15070		ם ם	0.00029	2610 17
61	20.99	4722		9	0.01000	026.04
01	24.07	4723		G	0.00472	920.94
		5885877			3.68762	1.29e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98827	Date	: 09/29/2009 07:51:21
Operator Sample Number	tchrom	Sample Name	: RSI0634-02
AutoSampler	: BUILT-IN	Rack/Vial	: 1/41
Instrument Name Instrument Serial #	: HP6890-06 : CN10520010	A/D mV Range	: B : 1000
Delay Time Sampling Rate	: 0.00 min : 5.0000 pts/s	End Time	: 29.45 min
Sample Volume	: 1.00000 ul	Area Reject	: 1500.000000
Data Acquisition Time	: 09/29/2009 03:08:27	Cycle	: 15

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46241.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46241.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46241.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46241.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46241.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	12433		В	0.01243	4528.01
2	9.42	1677		В	0.00168	614.05
3	10.54	15152	Tetrachloro-m-xylene (В	0.00221	4868.81
4	10.92	10143		В	0.01014	2787.09
5	11.91	9057		В	0.00906	2552.93
6	12.00	3233		V	0.00323	1000.04
7	12.16	3731		В	0.00373	645.77
8	12.31	688295	alpha-BHC (2C)	V	0.04648	217183.77
9	13.18	6196	,	В	0.00620	1769.47
10	13.26	12824	gamma-BHC (2C)	V	0.00231	2802.48
11	13.48	102075	beta-BHC (2C)	В	0.01684	29483.26
12	14.23	16819	delta-BHC (2C)	В	0.00323	5100.62
13	14.49	16813	()	В	0.01681	4794.67
14	14.66	3064		В	0.00306	846.52
15	14.94	3238		В	0.00324	713.81
16	15.35	7009		В	0.00701	2124.16

09/29/2009 07:51:21 Result: H:\TURBO6\6890-06\6-SEQ46\6b46241.rst

Peak #	Time [min]	Area [uV-sec]	Component Name		NG CONCENTRATION	Height [µV]
17	15.86	2972		В	0.00297	627.37
18	16.06	2855		В	0.00285	992.77
19	16.26	57716		B	0.05772	14383.99
20	16.55	2038	Hept. epoxide (2C) N	FВ	2.11e-04	546.70
21	16.77	3426		В	0.00343	1173.36
22	17.12	48127	(В	0.04813	7766.76
23	17.35	15869	alpha chlordane (2C)	V	0.00157	5113.51
24	17.77	19805		В	0.01980	4156.05
25	18.01	3065	Dieldrin (2C)	В	0.00118	905.18
26	18.52	2451		В	0.00245	780.53
27	18.62	1882		V	0.00188	546.05
29	18.96	3562	4,4'-DDD (2C)	V	0.00155	1134.78
30	19.23	2338	Endosulfan II (2C)	В	2.51e-04	675.22
31	19.67	3843	4,4'-DDT (2C)	В	0.00286	1008.69
32	19.90	5255	Endrin aldehyde (2C)	В	-1.2e-04	1574.95
33	20.04	3839		V	0.00384	1107.87
34	20.28	25785		В	0.02578	2957.20
35	20.44	102120		V	0.10212	23971.72
36	20.84	122455		В	0.12246	10265.28
37	21.15	20623	Methoxychlor (2C)	В	0.00565	5283.38
38	21.47	8801		В	0.00880	2061.15
39	21.63	2521		/ B	0.00252	802.40
40	21.77	35279	Endrin ketone (2C) 🐧	/ V	0.00512	7532.63
41	22.24	32880	V	В	0.03288	6820.94
43	23.10	208791	•	В	0.20879	38929.88
44	23.49	31376		В	0.03138	2419.85
45	23.75	10276		V	0.01028	1858.58
46	24.13	10115		В	0.01012	1448.69
47	24.47	4702		В	0.00470	1029.90
48	24.62	1867		В	0.00187	543.20
49	24.78	46788		V	0.04679	5933.57
50	25.25	49481	Decachlorobiphenyl (2C	В	0.00373	7443.10
51	25.65	28901		В	0.02890	4370.11
52	27.34	35951		В	0.03595	4177.81
53	27.56	24948		V	0.02495	2251.66
		1894462			1.01100	454410.34

31/275

Form 1 ORGANIC ANALYSIS DATA SHEET

CGS-DS1-091709

8081A

Laboratory:	TestAmerica Buffalo				SDG:			
Client: Olin Chlor Alkali Products - Cleveland,		<u>d, TN</u>	<u>'N</u> Project:		Olin - Charles Gibson site - NY3A9025AE03759			
Matrix:	<u>Solid</u>	Laborato	ry ID: <u>R</u>	RSI0634	-03	File ID:	<u>6A46242</u>	
Sampled:	09/17/09 12:00	Prepared	: <u>0</u>	9/21/09	08:00	Analyzed:	<u>09/29/09 03:44</u>	
Solids:	27.16	Preparati	on: <u>3</u>	550B G	C	Initial/Final:	<u>30.54 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>RI92925</u>		Calibration:	<u>R9I2904</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CONC	C. (ug/kg)	Q
319-84-6	alpha-BHC			10	210		D	
319-85-7	beta-BHC				10	73		D
319-86-8	delta-BHC				10	32		ற
58-89-9	gamma-BHC (Lindane)				10	60		UD
SYSTEM MONITORING COMPOUND			ADDED (u	ıg/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q
Decachlorobiphenyl			24.1		27.0	112	42 - 146	
Tetrachloro-m-xylene			24.1		28.9	120	37 - 135	

* Values outside of QC limits
| Software Version | : 6.2.1.0.104:0104 | Date : 09/29/2009 07:51:25 | |
|-----------------------|-------------------------|----------------------------|---|
| Reprocess Number | : buf2048: 98828 | | |
| Operator | : tchrom | Sample Name : RSI0634-03 | |
| Sample Number | : | Study : | |
| AutoSampler | : BUILT-IN | Rack/Vial : 1/42 | |
| Instrument Name | : HP6890-06 | Channel : A | |
| Instrument Serial # | : CN10520010 | A/D mV Range : 1000 | |
| Delay Time | : 0.00 min | End Time : 29.47 min | 4 |
| Sampling Rate | : 5.0000 pts/s | | |
| Sample Volume | : 1.00000 ul | Area Reject : 1500.000000 | |
| Sample Amount | : 1.0000 | Dilution Factor : 10.00 | |
| Data Acquisition Time | e : 09/29/2009 03:44:16 | Cycle : 16 | |

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46242.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46242.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46242.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46242.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46242.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	7.42	7594		В	0.00759	658.36
2	7.67	8621		В	0.00862	2087.21
3	7.77	31541		V	0.03154	9904.84
4	8.82	10993		В	0.01099	4281.42
5	9.09	15673		В	0.01567	5046.37
6	9.19	38373	Tetrachloro-m-xylene	V	0.00240	12222.87
7	9.94	2367	•	В	0.00237	1003.43
8	10.28	10052		В	0.01005	3423.53
9	10.76	661004	alpha-BHC	В	0.01782	217474.41
10	11.19	5071	•	В	0.00507	623.99
11	11.61	6351-	-gamma-BHC	В	0.00134	1729.11
12	11.86	93643	beta-BHC	В	0.00602	28287.71
13	12.30	25055	delta-BHC	В	0.00262	7801.63
14	12.47	18355		В	0.01836	5902.74
15	12.79	4247	Heptachlor NU	В	3.62e-04	1298.85
17	13.39	1845	- 1	В	0 00184	1059.03

09/29/2009 07:51:25 Result: H:\TURBO6\6890-06\6-SEQ46\6A46242.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	13.54	6088	Aldrin	NFE	B 0.00110	780.62
19	14.17	11758	r	í B	B 0.01176	2472.38
20	14.45	26631		B	0.02663	1822.47
21	14.62	95269		V	0.09527	22863.43
22	14.95	1713	Hept. epoxide	В	1.08e-05	585.55
23	15.27	26685	gamma chlordane) B	8 0.00168	7690.91
24	15.43	48087	-	E	0.04809	12267.63
25	15.71	13380) E	0.01338	3020.97
26	15.79	13861	4,4'-DDE	V	0.00169	3273.11
27	15.91	13305	Endosulfan I	V	5.58e-04	3435.48
28	16.02	6939		V	0.00694	1754.80
29	16.20	22127		E	0.02213	5727.69
30	16.43	3181	Dieldrin	B	0.00104	1154.01
31	16.64	2057		B	0.00206	756.71
32	16.95	35330	Endrin	B	0.00262	12639.43
33	17.10	3724	4,4'-DDD	B	0.00121	1378.31
34	17.72	20650	4,4'-DDT	B	0.00344	3226.79
35	18.15	3907		B	0.00391	1153.15
36	18.25	13424	Endrin aldehyde	, V	0.00101	3831.68
37	18.58	157315		ЫВ	0.15732	40239.07
38	18.89	9069) в	0.00907	2283.26
39	19.05	79656		\ V	0.07966	15305.46
40	19.38	51867		\ В	0.05187	9519.68
41	19.59	10092			0.01009	2543.15
42	19.75	5055	N	/ B	0.00505	1943.19
43	19.83	6670	Endrin ketone	J/ V	0.00229	1790.16
44	20.16	115677	,	V B	0.11568	20282.71
45	20.73	365667		В	0.36567	83784.25
46	21.08	33688		B	0.03369	9423.79
47	21.62	30826		B	0.03083	4694.45
48	21.70	37903		V	0.03790	5784.07
49	21.99	35306		B	0.03531	7628.45
50	22.09	69518	Decachlorobiphen	yl V	0.00224	13379.18
51	22.47	1094443		V	1.09444	93263.38
52	22.81	2103		B	0.00210	678.73
53	23.10	28483		B	0.02848	4312.40
54	23.69	36331		В	0.03633	6535.12
55	24.88	8420		B	0.00842	1170.43
		3486989			2.49362	717201.59

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98829	Date	: 09/29/2009 07:51:28
Operator	: tchrom	Sample Name	: RSI0634-03
Sample Number	• •	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/42
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 03:44:16	Cycle	: 16

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46242.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46242.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46242.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46242.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46242.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	15606		В	0.01561	5639.54
2	9.41	5341		В	0.00534	1868.73
3	9.65	1767		В	0.00177	272.40
4	10.09	1917		В	0.00192	634.62
5	10.54	15196	Tetrachloro-m-xylene (В	0.00221	4823.49
6	10.91	7398	2	В	0.00740	2125.38
7	11.91	6517		В	0.00652	1798.63
8	12.00	3073		V	0.00307	966.06
9	12.13	3554		В	0.00355	623.52
10	12.30	277392	alpha-BHC (2C)	V	0.02003	87098.97
11	13.18	1855	gamma-BHC (2C) RT	В	0.00154	677.11
12	13.48	38802	beta-BHC (2C)	В	0.00576	11535.04
13	14.22	10454	delta-BHC (2C)	В	0.00279	3137.68
14	14.49	8732	, , , , , , , , , , , , , , , , , , ,	В	0.00873	2567.30
16	14.93	5114		В	0.00511	853.41
17	15.35	9405		В	0.00941	2879.18

09/29/2009 07:51:28 Result: H:\TURBO6\6890-06\6-SEQ46\6b46242.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	16.05	2185		В	0.00218	785.63
19	16.26	37802		В	0.03780	9610.20
20	16.77	3223		В	0.00322	1067.43
21	17.02	2926	gamma chlordane (2C)	MB	7.15e-04	1097.41
22	17.12	33108		V"V	0.03311	5503.98
23	17.34	16028	alpha chlordane (2C)	V	0.00158	5337.69
24	17.77	13178		/ В	0.01318	2762.76
25	18.01	2006	Dieldrin (2C)	В	0.00108	612.39
28	18.96	3053	4,4'-DDD (2C)	V	0.00149	941.69
29	19.22	1507	Endosulfan II (2C)	В	1.59e-04	498.49
30	19.67	1679	4,4'-DDT (2C)	В	0.00259	487.25
31	19.90	3947	Endrin aldehyde (2C)	В	-3.4e-04	1232.14
32	20.04	2013	1	В	0.00201	620.57
33	20.28	25454		В	0.02545	2808.17
34	20.44	77845		V	0.07785	16892.86
35	20.77	90193		V	0.09019	10261.53
36	20.84	66947		V	0.06695	10789.00
37	20.96	29063		V	0.02906	4344.31
38	21.16	20357	Methoxychlor (2C)	V	0.00558	4295.14
39	21.47	10045	ł	В	0.01004	1855.53
40	21.63	3651	Ý	V	0.00365	829.50
41	21.77	19752	Endrin ketone (2C)	V	0.00337	3779.17
42	22.22	406611		V	0.40661	23325.42
43	23.10	147718		В	0.14772	27942.57
44	23.50	4993		В	0.00499	389.90
45	23.76	14569		В	0.01457	2865.63
48	24.78	28877		В	0.02888	3930.06
49	25.25	46630	Decachlorobiphenyl (20	; В	0.00339	6371.58
50	25.65	20279		В	0.02028	3030.91
51	27.34	13756		В	0.01376	1957.14
		1561519			1,15189	283727.11

CGS-DSD1-091709

ORGANIC ANALYSIS DATA SHEET

8081A

Laboratory:	TestAmerica Buffalo				SDG:			
Client:	<u>Olin Chlor Alkal</u>	i Products - Clevel	and, TN		Project: <u>Olin - Charles Gibson site - NY3A9</u>			25AE03759
Matrix:	<u>Solid</u>	Labora	tory ID:	<u>RSI0634</u>	<u>1-04</u>	File ID:	<u>6A46243</u>	
Sampled:	<u>09/17/09 12:20</u>	Prepare	ed:	<u>09/21/09</u>	<u>0 08:00</u>	Analyzed:	<u>09/29/09 04:20</u>	
Solids:	<u>37.92</u>	Prepara	ation:	<u>3550B (</u>	<u>GC</u>	Initial/Final:	<u>30.44 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>R192925</u>		Calibration:	<u>R9I2904</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CON	C. (ug/kg)	Q
319-84-6	alpha-BHC				10		93	D
319-85-7	beta-BHC				10		170	D
319-86-8	delta-BHC				10	24		л
58-89-9	gamma-BHC (Li	ndane)			10		43	UD
SYSTEM MONI	TORING COMPO	UND	ADDED) (ug/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q
Decachlorobiphenyl		17	7.3	21.2	122	42 - 146		
Tetrachloro-m-xy	ylene		17	7.3	20.7	119	37 - 135	

* Values outside of QC limits

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98830	Date	: 09/29/2009 07:51:31
Operator	: tchrom	Sample Name	: RSI0634-04
Sample Number	:	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/43
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 04:20:15	Cycle	: 17

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46243.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46243.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46243.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46243.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46243.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.42	12979		в	0.01298	1535.07
2	7.51	7348		V	0.00735	2375.95
3	7.66	20391		В	0.02039	5250.68
4	7.77	44240		V	0.04424	13094.03
5	8.82	6588		В	0.00659	2693.10
6	9.09	21584		В	0.02158	6880.27
7	9.19	38121	Tetrachloro-m-xylene	V	0.00239	12040.67
8	9.94	11073	•	В	0.01107	2607.57
9	10.28	12934		В	0.01293	4283.49
10	10.76	382604	alpha-BHC	В	0.01075	125590.32
11	11.19	4617	•	В	0.00462	608.27
12	11.33	10200		В	0.01020	2994.27
13	11.61	4618	gamma-BHC PT	В	0.00129	1441.05
14	11.86	285288	beta-BHC	В	0.01956	88134.90
15	12.30	31136	delta-BHC	В	0.00278	9877.96
16	12.47	34835	*	В	0.03484	11146.03

09/29/2009 07:51:31 Result: H:\TURBO6\6890-06\6-SEQ46\6A46243.rst

Peak #	Time		Component	BL		Height
#	frond	[uv-sec]	INdifie		CONCENTRATION	[h ʌ]
17	12.79	17511	Heptachlor M	В	7.58e-04	3230.18
19	13.39	2292	instruction with	B	0.00229	1331.06
20	13.53	5955	Aldrin	B	0.00110	952.15
21	14.17	9320		B	0.00932	2443.66
22	14.32	5949	\	B	0.00595	911.09
23	14.62	112759	\backslash	B	0.11276	28345.77
24	15.00	11048	Hept, epoxide	B	3.13e-04	1497.21
25	15.27	28517	gamma chlordane	B	0.00174	8145.06
26	15.43	62633	3	В	0.06263	15428.16
27	15.72	61527	4.4'-DDE	В	0.00320	20981.57
28	15.91	7554	Endosulfan I	В	3.52e-04	2932.50
29	16.02	6088		В	0.00609	1923.39
30	16.20	28285		B	0.02829	7314.62
31	16.43	3871	Dieldrin	\ B	0.00106	1364.78
32	16.64	2559		B	0.00256	946.65
33	16.95	36449	Endrin	B	0.00266	12993.26
35	17.71	25048	4,4'-DDT	В	0.00362	3602.15
36	18.15	6524		B	0.00652	1828.65
37	18.25	17863	Endrin aldehyde	/V	0.00127	4800.87
38	18.58	194187	-	B	0.19419	49962.53
39	18.89	10608		В	0.01061	2809.14
40	19.05	105160		V	0.10516	20899.75
41	19.38	64575		B	0.06458	11820.08
42	19.59	13045		V	0.01305	3270.05
43	19.75	12698		[V	0.01270	3140.06
44	19.82	17587	Endrin ketone	V	0.00273	3746.85
45	20.16	136102		B	0.13610	24540.32
46	20.73	460137	ľ	B	0.46014	105343.70
47	21.08	20181		[∕B	0.02018	5599.93
48	21.59	62825		'В	0.06282	11046.72
49	21.99	50158		В	0.05016	10099.65
50	22.09	74093	Decachlorobiphenyl	V	0.00245	14132.30
51	22.47	760812		V	0.76081	62635.54
52	22.63	64357		V	0.06436	14238.94
53	23.10	27066		В	0.02707	4658.33
54	23.69	53261		В	0.05326	8659.79
55	23.98	7754		В	0.00775	1670.16
56	24.88	9371		В	0.00937	1158.59
		3524284			2.53352	770958.82

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98831	Date	: 09/29/2009 07:51:34
Operator	: tchrom	Sample Name	: RSI0634-04
Sample Number	:	Study	•
AutoSampler	: BUILT-IN	Rack/Vial	: 1/43
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/29/2009 04:20:15	Cycle	: 17

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46243.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46243.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46243.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46243.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46243.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	18691		В	0.01869	6707.12
2	9.42	3282		В	0.00328	1099.22
3	9.64	1618		В	0.00162	322.09
4	10.09	3403		В	0.00340	1116.76
5	10.54	15088	Tetrachloro-m-xylene (В	0.00220	4795.47
6	10.91	10258	2 (В	0.01026	2977.56
7	11.62	1949		В	0.00195	715.94
8	11.77	4537		В	0.00454	1281.06
9	11.90	9872		V	0.00987	2447.44
10	12.00	3752		V	0.00375	1126.98
11	12.15	3392		В	0.00339	558.75
12	12.30	163742	alpha-BHC (2C)	V	0.01272	51032.18
13	13.17	2752	gamma-BHC (2C)	В	0.00160	926.62
14	13.48	126824	beta-BHC (2C)	В	0.02117	36875.15
15	13.91	2878	5	В	0.00288	122.06
16	14.23	13268	delta-BHC (2C)	В	0.00298	4043.62

09/29/2009 07:51:34 Result: H:\TURBO6\6890-06\6-SEQ46\6b46243.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	14.49	17461		B	0.01746	5029.50
19	14.93	6920		В	0.00692	1173.84
20	15.35	10647		В	0.01065	3159.05
21	15.86	4184		В	0.00418	888.02
22	16.06	28624		В	0.02862	9412.40
23	16.27	51187		В	0.05119	12191.54
24	16.77	4502		В	0.00450	1391.78
25	17.12	40628		B	0.04063	6504.42
26	17.34	16496	alpha chlordane (2C)	1/PAV	0.00163	5596.97
27	17.77	15681		B	0.01568	3364.91
28	18.00	2760	Dieldrin (2C)	/ B	0.00115	804.94
29	18.42	9186		В	0.00919	877.78
30	18.51	2311		V	0.00231	739.16
31	18.96	2261	4,4'-DDD (2C)	В	0.00140	802.52
32	19.22	3171	Endosulfan II (2C)	B	3.45e-04	917.59
33	19.67	2729	4,4'-DDT (2C)	В	0.00272	697.05
34	19.90	6307	Endrin aldehyde (2C)	B	5.34e-05	1753.86
35	20.04	4430		v	0.00443	1182.49
36	20.29	17125		В	0.01713	1913.02
37	20.44	94301		V	0.09430	21007.34
38	20.60	5529	Endo. Sulfate (2C)	E	0.00270	1693.63
39	20.84	115253		V	0.11525	9874.75
40	20.97	26281		V	0.02628	3921.85
41	21.15	24549	Methoxychlor (2C)	v	0.00671	5337.36
42	21.47	11414		В	0.01141	2094.14
43	21.63	5775		_ / V	0.00578	1225.10
44	21.77	32577	Endrin ketone (2C)	/ V	0.00482	6063.71
45	22.23	306166	V	r V	0.30617	18920.64
46	23.10	180556		В	0.18056	34023.12
47	23.49	2375		В	0.00238	254.73
48	23.75	9202		В	0.00920	1786.43
49	24.12	1695		В	0.00170	317.20
50	24.46	3833		В	0.00383	865.18
51	24.78	36753		В	0.03675	4940.77
52	25.25	52070	Decachlorobiphenyl (2	СВ	0.00403	7134.79
53	25.64	25059		В	0.02506	3747.87
54	27.34	21161		В	0.02116	2892.96
56	28.40	1582		В	0.00158	146.16
57	29.37	2644		В	0.00264	219.20
		1590690			1.18680	301015.79

41/275

Form 6

INITIAL CALIBRATION DATA

Bit Bit <th colspan="5">Laboratory: <u>TestAmerica Buffalo</u></th> <th colspan="5">SDG:</th>	Laboratory: <u>TestAmerica Buffalo</u>					SDG:								
Burner	Client:	<u>Olin Chlo</u>	or Alka	li Products -	Clevela	ind, TN		Project:		<u>Olin -</u>	Charle	s Gibson site	<u>- NY3.</u>	A9025AE03
neuron neuron<	Calibration:	R9H1803	3					Instrume	ent:	HP68	90-6			
Level O1 Level O2 Level O3 Level O3 Level O3 Level O3 Level O3 Level O3 Reb			-					Calibrati	an Date		<u>/00 10.(</u>			
<table-container> Processe Processe</table-container>								Calibrati	on Date	s: <u>07/25</u>	/09/10:0	<u>15</u>		
Compoundnynynynynynynynynynynyny4A*DDDii			L	evel 01	L	evel 02	L	evel 03	Level 04		L	evel 05	Level 06	
44-DDD1111111122244-DDD11	Compound		ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF
44-0D(A·D)(A)<	4,4'-DDD									1		.	0.15	2.247233E+07
44-0DE1.4. <th< td=""><td>4,4'-DDD [2C]</td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td>0.15</td><td>7274867</td></th<>	4,4'-DDD [2C]												0.15	7274867
4.4-DDP(1)(1)(1)(1)(1)(1)(1)(1)(1)(1)(2)(2)4.4-DDP(1)	4,4'-DDE								1				0.15	2.656103E+07
44-DDTimimimimimimimimimimim4A-DDT (2)im <td>4,4'-DDE [2C]</td> <td></td> <td></td> <td>1</td> <td></td> <td></td> <td></td> <td></td> <td>1</td> <td></td> <td></td> <td></td> <td>0.15</td> <td>9062533</td>	4,4'-DDE [2C]			1					1				0.15	9062533
44-DIC[C]101101101101101101101101101101Aldin101101101101101101101101101101101alpha-BHC101101101101101101101101101101101101alpha-BHC10110	4,4'-DDT				1					1	<u> </u>		0.15	2.353113E+07
AddmAddmCiC	4,4'-DDT [2C]												0.15	6604504
Aldin [2C] In	Aldrin										<u> </u>		0.15	2.540948E+07
alpha-BHC alpha-BHC (2C)ininininininininininininalpha-BHC alpha-Chordame [2C]in<	Aldrin [2C]			<u> </u>			<u> </u>				<u> </u>		0.15	1.086657E+07
alpha-BHC [2C]in<in<in<<	alpha-BHC					ļ	<u> </u>		1			<u> </u>	0.15	3.138012E+07
alpha-ChloradaeIn </td <td>alpha-BHC [2C]</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>·</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td>0.15</td> <td>1.524203E+07</td>	alpha-BHC [2C]						·						0.15	1.524203E+07
apha-Chordane [2C] i.e. i.e. <td>alpha-Chlordane</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td><u> </u></td> <td></td> <td>0.15</td> <td>2.489279E+07</td>	alpha-Chlordane										<u> </u>		0.15	2.489279E+07
beta-BHCii </td <td>alpha-Chlordane [2C]</td> <td></td> <td></td> <td></td> <td></td> <td></td> <td></td> <td> </td> <td><u> </u></td> <td></td> <td></td> <td></td> <td>0.15</td> <td>9171053</td>	alpha-Chlordane [2C]								<u> </u>				0.15	9171053
base-BHC [2C] i <	beta-BHC			<u> </u>			<u> </u>	· · · · ·				1	0.15	1 139781E+07
base breach	beta-BHC [2C]			l					<u> </u>				0.15	5422020
Decentionsplantly 0.0 2.000000000000000000000000000000000000	Decachlorohinhenul		0.1	2 2220425+02	0.076	2 2202215:07	0.05	2 21215 (02		2.45(2845)07	0.005	D 27062E+02	0.15	5432323
Deck monomenty (EQ) 0.11 79/92 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 0005 799/90 799	Decachlorobiphenyl [2	PC1	0.1	2.22/0422+0/	0.075	2.328/31040/	0.05	2.21218+07	0.01	2.430264E+07	0.005	2.37932E+07	<u> </u>	· · · · · · · · · · · · · · · · · · ·
actus BHC i	Decacillorooipnenyi [2		0.1	7191925	0.075	7389729	0.05	7399140	0.01	8690760	0.005	8993840		
delta Phi [2C] i	denta-BHC												0.15	2.946435E+07
Diekirin Image Image <thimage< th=""> Image Image <</thimage<>	Delta-BHC [2C]			ļ	ļ								0.15	1.3229E+07
Deckmin [AC] I <t< td=""><td>Dielarin</td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td>ļ</td><td> </td><td><u> </u></td><td></td><td>0.15</td><td>2.830859E+07</td></t<>	Dielarin								ļ		<u> </u>		0.15	2.830859E+07
Endosultan I	Dieldrin [2C]			·			ļ						0.15	9453413
Endosulfan I [2C] Image: Marcine Marci	Endosultan I				_				ļ				0.15	2.433352E+07
Endosulfan II <	Endosulfan I [2C]												0.15	8616140
Endosulfan II [2C] I	Endosulfan II												0.15	2.360599E+07
Endosulfan sulfate Image: Marking Mark	Endosulfan II [2C]												0.15	7630460
Endosulfan sulfate [2C]II <t< td=""><td>Endosulfan sulfate</td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td>0.15</td><td>2.164311E+07</td></t<>	Endosulfan sulfate												0.15	2.164311E+07
EndrinImage: start of the start	Endosulfan sulfate [2C]]						0.15	7065894
Endrin [2C]Image: second s	Endrin												0.15	2.498895E+07
Endrin aldehydeImage: sector of the sector of t	Endrin [2C]												0.15	7605587
Endrin aldehyde [2C]Image: sector of the sector	Endrin aldehyde												0.15	1.760603E+07
Endrin ketoneImage: second	Endrin aldehyde [2C]												0.15	5554784
Endrin ketone [2C]III<	Endrin ketone												0.15	2.520435E+07
gamma-BHC (Lindane) Image: Second Secon	Endrin ketone [2C]												0.15	8157214
gamma-BHC (Lindane) [2C] Image: Constraint of the symbol of	gamma-BHC (Lindane	;)											0.15	2.860553E+07
gamma-Chlordane Image: state in the state i	gamma-BHC (Lindane	e) [2C]										· · · ·	0.15	1.345126E+07
gamma-Chlordane [2C] I	gamma-Chlordane												0.15	2.72336E+07
Heptachlor Image: Marking Mark	gamma-Chlordane [2C	;]											0.15	9720073
Heptachlor [2C] Image: Constraint of the system of the syste	Heptachlor												0.15	2.632172E+07
Heptachlor epoxide Image: Constraint of the system Image: Consystem Image: Constraint of the system	Heptachlor [2C]										<u> </u>		0.15	1.186545E+07
Heptachlor epoxide [2C] Image: Constraint of the second secon	Heptachlor epoxide												0.15	2.537119E+07
Methoxychlor Image: Constraint of the second s	Heptachlor epoxide [20			├							<u> </u>	<u> </u>	0.15	9762733
Methoxychior [2C] Image: Constraint of the second sec	Methoxychlor								<u> </u>				0.15	1.171084F+07
Tetrachloro-m-xylene 0.1 1.752098E+07 0.075 1.610821E+07 0.05 1.615162E+07 0.01 1.521668E+07 0.005 1.496592E+07 0 Tetrachloro-m-xylene 0.1 1.012139E+07 0.075 9238272 0.05 9281484 0.01 9583810 0.005 9486640	Methoxychlor [2C]												0.15	3362157
Tetrachloro-m-xylene [2C] 0.1 1.010216107 0.05 1.010216107 0.05 1.01362E107 0.01 1.52100E207 0.005 1.490592E107	Tetrachloro-m-vvlene		0.1	1.7520088407	0.075	1.6108215-07	0.05	16151678-07	0.01	1 5716695±07	0.005	1 4965025+07		
0.0 1.01 1.01 1.01 1.000	Tetrachloro-m-xvlene [[2C]	0.1	1.012139E+07	0.075	9738777	0.05	9781484	0.01	9583810	0.005	9486640		

INITIAL CALIBRATION DATA (Continued)

		Lovel 07	Lovel 08	Louol 00	Level 10	Lovel 11	Lovel 12	
Calibration Date: 07/23/09 10:05								
Calibration:	<u>R9H1803</u>			Instrume	nt: <u>HP68</u>	<u>90-6</u>		
Client:	Olin Chlo	or Alkali Products -	Cleveland, TN	Project:	<u>Olin</u> ·	- Charles Gibson sit	<u>e - NY3A9025AE03'</u>	
Laboratory:	TestAme	rica Buffalo		SDG:				

	L	Level 07		Level 08		evel 09	L	evel 10	Le	evel 11	L	evel 12
Compound	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF
4,4'-DDD	0.1	2.168968E+07	0.05	2.156712E+07	0.01	1.839035E+07	0.005	1.81645E+07				
4,4'-DDD [2C]	0.1	6934292	0.05	6921504	0.01	6266309	0,005	6402734				
4,4'-DDE	0.1	2.535969E+07	0.05	2.500088E+07	0.01	2.060628E+07	0.005	2.0436E+07				
4,4'-DDE [2C]	0.1	8619720	0.05	8545540	0.01	7601200	0.005	7822780				
4,4'-DDT	0.1	2.272436E+07	0.05	2.253244E+07	0.01	1.83517E+07	0.005	1.802112E+07				
4,4'-DDT [2C]	0.1	6366408	0.05	6296356	0.01	5303650	0.005	5233760		•		
Aldrin	0.1	2.421195E+07	0.05	2.389178E+07	0.01	1.931586E+07	0.005	1.90514E+07				
Aldrin [2C]	0.1	1.032205E+07	0.05	1.019713E+07	0.01	8969080	0.005	9278140				
alpha-BHC	0.1	2.951163E+07	0.05	2.845386E+07	0.01	2.170381E+07	0.005	2.087144E+07				
alpha-BHC [2C]	0.1	1.427292E+07	0.05	1.385803E+07	0.01	1.118012E+07	0.005	1.101172E+07				
alpha-Chlordane	0.1	2.378313E+07	0.05	2.39888E+07	0.01	2.130226E+07	0.005	2.214108E+07				
alpha-Chlordane [2C]	0.1	8826526	0.05	9005336	0.01	9311152	0.005	1.051486E+07				
beta-BHC	0.1	1.086216E+07	0.05	1.103278E+07	0.01	1.036334E+07	0.005	1.068322E+07				
beta-BHC [2C]	0.1	5255053	0.05	5491150	0.01	5531920	0.005	5785520				
delta-BHC	0.1	2.754822E+07	0.05	2.662916E+07	0.01	2.05191E+07	0.005	1.98651E+07				
delta-BHC [2C]	0.1	1.238175E+07	0.05	1.218441E+07	0.01	1.015513E+07	0.005	1.02742E+07				
Dieldrin	0.1	2.713924E+07	0.05	2.693052E+07	0.01	2.322494E+07	0.005	2.331834E+07				
Dieldrin [2C]	0.1	9028876	0.05	9022780	0.01	8275660	0.005	8562360				
Endosulfan I	0.1	2.353209E+07	0.05	2.380174E+07	0.01	2.143969E+07	0.005	2.20314E+07				
Endosulfan I [2C]	0.1	8317141	0.05	8482628	0.01	8198468	0.005	8629376				
Endosulfan II	0.1	2.26842E+07	0.05	2.331696E+07	0.01	2.117094E+07	0.005	2.197706E+07				
Endosulfan II [2C]	0.1	7390036	0.05	7679686	0.01	7493460	0.005	7930840				
Endosulfan sulfate	0.1	2.072924E+07	0.05	2.147146E+07	0.01	1.971092E+07	0.005	2.057732E+07				
Endosulfan sulfate [2C]	0.1	6856330	0.05	7307560	0.01	7280250	0.005	7758360				
Endrin	0.1	2.424614E+07	0.05	2.404716E+07	0.01	2.264809E+07	0.005	2.403058E+07				
Endrin [2C]	0.1	7319696	0.05	7311648	0.01	6836871	0.005	6965468				
Endrin aldehyde	0.1	1.691286E+07	0.05	1.780957E+07	0.01	1.623091E+07	0.005	1.715176E+07				
Endrin aldehyde [2C]	0.1	5376208	0.05	5761344	0.01	5763500	0.005	6160000				
Endrin ketone	0.1	2.425007E+07	0.05	2.474564E+07	0.01	2.195196E+07	0.005	2.242924E+07				
Endrin ketone [2C]	0.1	7924476	0.05	8281502	0.01	7810021	0.005	8137040				
gamma-BHC (Lindane)	0.1	2.698992E+07	0.05	2.626372E+07	0.01	2.089465E+07	0.005	2.042404E+07				
gamma-BHC (Lindane) [2C]	0.1	1.271798E+07	0.05	1.25653E+07	0.01	1.064662E+07	0.005	1.080584E+07				
gamma-Chlordane	0.1	2.59836E+07	0.05	2.583742E+07	0.01	2.288938E+07	0.005	2.365148E+07				
gamma-Chlordane [2C]	0.1	9301344	0.05	9372676	0.01	9070590	0.005	9744520				
Heptachlor	0.1	2.539545E+07	0.05	2.534458E+07	0.01	2.152437E+07	0.005	2.15986E+07				
Heptachlor [2C]	0.1	1.137805E+07	0.05	1.13983E+07	0.01	1.035612E+07	0.005	1.082652E+07				
Heptachlor epoxide	0.1	2.455175E+07	0.05	2.479796E+07	0.01	2.1992E+07	0.005	2.21962E+07				
Heptachlor epoxide [2C]	0.1	9417434	0.05	9518412	0.01	9116520	0.005	9611320				
Methoxychlor	0.1	1.161118E+07	0.05	1.200439E+07	0.01	1.111994E+07	0.005	1.123854E+07				
Methoxychlor [2C]	0.1	3375054	0.05	3551314	0.01	3242100	0.005	3227200				

INITIAL CALIBRATION DATA (Continued)

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Laboratory:	<u>TestAmerica Buffa</u>	<u>lo</u>		SD	G:				
Client:	<u>Olin Chlor Alkali F</u>	Products - Cleve	eland, TN	Pro	ject:	<u>Olin - Cha</u>	rles Gibson site	<u>- NY3A9025AE</u>	03759
Calibration:	<u>R9H1803</u>			Inst	rument:	<u>HP6890-6</u>			
				0-1	il	07/22/00 1	0.05		
				Cal	ibration Date:	0//23/09 1	0:05		
Compound		Mean RF	RF RSD	Mean RT	RT RSD	Linear r ²	Quad COD	LIMIT	Q
4,4'-DDD		2.04568E+07	9.880035	17.58529	1.697931E-02	0.99979			
4,4'-DDD [2C]		6759941	6.156987	19.49521	2.119975E-02	0.99956			
4,4'-DDE		2.359278E+07	12.13739	16.25639	1.700435E-02	0.99962			
4,4'-DDE [2C]		8330355	7.241039	18.21338	2.370574E-02	0.99951			
4,4'-DDT		2.103215E+07	12.49175	18.19676	1.343948E-02	0.99980			
4,4'-DDT [2C]		5960936	10.78072	20.1964	1.784081E-02	0.99977			
Aldrin		2.237609E+07	13.27387	14.02588	3.264287E-02	0.99960			
Aldrin [2C]		9926595	7.884857	15.74869	2.079922E-02	0.99948			
alpha-BHC		2.638417E+07	18.10964	11.22486	4.305247E-02	0.99930			
alpha-BHC [2C]		1.311296E+07	14.56188	12.81126	3.369039E-02	0.99918			
alpha-Chlordane		2.322161E+07	6.295268	16.07044	3.218475E-03	0.99961			
alpha-Chlordane [20	C]	9365786	7.126577	17.8786	1.257882E-02	0.99965			
beta-BHC		1.086786E+07	3.553908	12.32182	3.697631E-02	0.99954			
beta-BHC [2C]		5499314	3.487293	13.98707	3.274282E-02	0.99971			
Decachlorobiphenyl	1	2.320735E+07	4.440189	22.68772	1.479249E-02	0.99935			
Decachlorobiphenyl	1 [2C]	7933079	10.60081	26.11406	3.233055E-02	0.99984			
delta-BHC		2.480519E+07	17.49535	12.76784	3.739646E-02	0.99916			
delta-BHC [2C]		1.16449E+07	11.713	14.74554	3.009804E-02	0.99918			
Dieldrin		2.578433E+07	9.127046	16.92149	2.941452E-02	0.99968			
Dieldrin [2C]		8868618	5.157413	18.64202	1.647727E-02	0.99959			
Endosulfan I		2.302769E+07	5.350876	16.37939	2.160991E-02	0.99979			
Endosulfan I [2C]		8448751	2.22816	18.02188	2.200692E-02	0.99974			
Endosulfan II		2.255103E+07	4.407369	17.937	2.067989E-02	0.99968			
Endosulfan II [2C]		7624897	2.696299	19.7883	2.201632E-02	0.99975			
Endosulfan sulfate		2.082641E+07	3.718438	19.81559	1.987521E-02	0.99961			
Endosulfan sulfate [2C]	7253679	4.631506	21.13427	1.490868E-02	0.99967			
Endrin		2.399218E+07	3.530003	17.43977	2.580875E-02	0.99980			
Endrin [2C]		7207854	4.264078	19.33217	1.633863E-02	0.99971			
Endrin aldehyde		1.714223E+07	3.623915	18.86289	1.271725E-02	0.99960			
Endrin aldehyde [20	C]	5723167	5.113908	20.52088	1.686058E-02	0.99961			
Endrin ketone		2.371625E+07	6.084135	20.42922	1.700597E-02	0.99972			
Endrin ketone [2C]		8062050	2.363925	22.44694	0.0157225	0.99977			
gamma-BHC (Linda	ane)	2.463557E+07	15.14537	12.07683	4.042926E-02	0.99938			
gamma-BHC (Linda	ane) [2C]	1.20374E+07	10.3359	13.7757	3.564005E-02	0.99941			
gamma-Chlordane		2.51191E+07	7.138167	15.76345	2.068199E-02	0.99958			
gamma-Chlordane [2C]	9441841	3.048426	17.544	0.0232897	0.99959			
Heptachlor		2.403694E+07	9.54012	13.29072	3.238207E-02	0.99978			
Heptachlor [2C]		1.116489E+07	5.221252	14.93749	2.670636E-02	0.99966			
Heptachlor epoxide		2.378182E+07	6.60485	15.47862	2.327685E-02	0.99981			
Heptachlor epoxide	[2C]	9485284	2.552927	17.10375	1.734474E-02	0.99973			ļ
Methoxychlor		1.153698E+07	3.116776	19.23775	2.164316E-02	0.99992			
Methoxychlor [2C]		3351565	3.890612	21.66569	1.537311E-02	0.99978			
Tetrachloro-m-xyler	ne	1.599268E+07	6.276828	9.636484	5.262336E-02	0.99817			
Tetrachloro-m-xyler	ne [2C]	9542319	3.707424	11.02601	0.0293783	0.99765			

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INITIAL CALIBRATION DATA

Laboratory:	TestAme	rica Bu	<u>iffalo</u>				SDG:						
Client:	<u>Olin Chle</u>	or Alka	li Products -	Clevela	ind, TN		Project:		<u>Olin -</u>	Charle	s Gibson site	<u>- NY3</u>	<u> 49025AE03'</u>
Calibration:	R912904						Instrume	nt:	HP68	90-6			
							Calibrati	on Date	e: <u>09/28</u>	/09 00:0	<u>90</u>		
		L	evel 01	L	evel 02	L	Level 03		evel 04	L	evel 05	L	evel 06
Compound		ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ui	RF
4,4'-DDD		0.15	2.338388E+07	0.1	2.367061E+07	0.05	2.316616E+07	0.01	1.986354E+07	0.005	1.94188E+07		
4,4'-DDD [2C]		0.15	8422420	0.1	8348680	0.05	8206096	0.01	7374160	0.005	7459721		
4,4'-DDE		0.15	3.105571E+07	0.1	3.122165E+07	0.05	3.051852E+07	0.01	2.600267E+07	0.005	2.564862E+07		
4,4'-DDE [2C]		0.15	1.080204E+07	0.1	1.0677E+07	0.05	1.046265E+07	0.01	9426680	0.005	9610760		
4,4'-DDT		0.15	2.440827E+07	0.1	2.369051E+07	0.05	2.271714E+07	0.01	1.854514E+07	0.005	1.770046E+07		
4,4'-DDT [2C]		0.15	7918940	0.1	7599594	0.05	7302550	0.01	6479320	0.005	6317800		
Aldrin		0.15	3.265655E+07	0.1	3.278402E+07	0.05	3.249588E+07	0.01	2.819759E+07	0.005	2.800768E+07		
Aldrin [2C]		0.15	1.229787E+07	0.1	1.231225E+07	0.05	1.205503E+07	0.01	1.075118E+07	0.005	1.091148E+07		
alpha-BHC		0.15	3.864773E+07	0.1	4.010065E+07	0.05	3.867142E+07	0.01	3.142344E+07	0.005	3.03854E+07		
alpha-BHC [2C]		0.15	1.537612E+07	0.1	1.513059E+07	0.05	1.452124E+07	0.01	1.200552E+07	0.005	1.183352E+07		
alpha-Chlordane		0.15	3.039283E+07	0.1	3.076903E+07	0.05	3.04763E+07	0.01	2.794036E+07	0.005	2.850292E+07		
alpha-Chlordane [20	C]	0.15	1.06857E+07	0.1	1.069717E+07	0.05	1.07186E+07	0.01	1.032234E+07	0.005	1.075731E+07		
beta-BHC		0.15	1.410899E+07	0.1	1.444747E+07	0.05	1.452133E+07	0.01	1.415674E+07	0.005	1.471012E+07		
beta-BHC [2C]		0.15	5718046	0.1	5820064	0.05	5945456	0.01	6022740	0.005	6288200		
Decachlorobipheny	1											0.1	2.250105E+07
Decachlorobipheny	1 [2C]							<u> </u>				0.1	8557185
delta-BHC		0.15	3.662848E+07	0.1	3.654457E+07	0.05	3.522454E+07	0.01	2.855362E+07	0.005	2.74564E+07		
delta-BHC [2C]		0.15	1.427003E+07	0.1	1.404673E+07	0.05	1.351333E+07	0.01	1.136683E+07	0.005	1.113774E+07		
Dieldrin		0.15	3.138933E+07	0.1	3.16393E+07	0.05	3.111638E+07	0.01	2.710447E+07	0.005	2.702184E+07		
Dieldrin [2C]		0.15	1.115007E+07	0.1	1.106901E+07	0.05	1.091491E+07	0.01	9997900	0.005	1.028246E+07		
Endosulfan I		0.15	2.775139E+07	0.1	2.817111E+07	0.05	2.81964E+07	0.01	2.604366E+07	0.005	2.649044E+07		
Endosulfan I [2C]		0.15	1.011625E+07	0.1	1.016266E+07	0.05	1.020055E+07	0.01	9791916	0.005	1.016429E+07	┟───┦	
Endosulfan II		0.15	2.488065E+07	01	2 510043E+07	0.05	2 493478E+07	0.01	2.342066E+07	0.005	2.348148E+07		
Endosulfan II [2C]		0.15	8034713	0.1	8043404	0.05	8064418	0.01	8773000	0.005	9093840		
Endosulfan sulfate		0.15	1943163E+07	0.1	1 860879E+07	0.05	1 77079E+07	0.01	1 511711E+07	0.005	1 40372E+07		
Endosulfan sulfate l	201	0.15	7100097	0.1	6000076	0.05	6758604	0.01	6195200	0.005	5848140		
Endrin		0.15	2 7202278+07	0.1	2 722051E+07	0.05	2 6661828407	0.01	2 197293E+07	0.005	2 251008E+07	┟╍╍──┦	
Endrin [2C]		0.15	0129572		0194142	0.05	2.0001822107	0.01	8180280	0.005	\$208960	[]	
Endrin aldehyde		0.15	1642246E+07	0.1	1 6604946-07	0.05	1 657505E±07	0.01	1.528704E+07	0.005	1 575264E+07	┣━━──┦	
Endrin aldehyde [2]		0.15	6000140	0.1	6056083	0.05	6101520	0.01	6300420	0.005	6737800	┟╾╺╍╾─┦	
Endrin ketone	~]	0.15	2 5009425-07	- 0.1	0030963	0.05	2 241786E±07	0.01	2 111154E+07	0.005	2 080012E+07	<u> </u>	
Endrin ketone [2C]		0.15	2.3006422707	0.1	2.3933462707	0.05	2.341/80240/	0.01	2.1111342107	0.005	9357190	┠────┦	
gamma_BHC (1 ind	ane)	0.15	2 5557(2)-07	0.1	000/000 0 ()(0)(0)	0.05	0049000	0.01	20629295-02	0.005	2 009002E+07	┟────┤	
gamma-BHC (Lind	ane) [2C]	0.15	3.333/6ETU/	0.1	1.400(25E+07	0.05	1.270644E+07	0.01	1 1959935+07	0.005	1 199252E+07	<u> </u>	
gamma-DHC (Entu		0.15	2.0720010-07	0.1	1.4096232707	0.05	1.3/0344E+07	0.01	2 980702E+07	0.005	2.01065+07	<u> </u>	
gamma-Chlordane	201	0.15	3.2/3201E+07	0.1	3.283498ETU/	0.03	3.230242E=07	0.01	2.889/92070	0.005	1.09552/E107	{l	
Hontachlar	20]	0.15	1.13489/E+0/	0.1	1.136915E+07	0.05	1.1285/2E+0/	0.01	1.052807E+07	0.005	1.065556E+07		
Hentachlor [2C]		0.15	1 222180E-07	0.1	3.301//0ETU/	0.05	1 2006/40107	0.01	1 2020295-02	0.005	1 2201548-07	<u> </u>	
Hentechler anorida		0.15	1.322189E+07	0.1	1.3252296+07	0.05	1.308566E+07	0.01	1.203028E+07	0.005	1.429130E+07		
Hentachlor mori -	1101	0.15	3.068174E+07	0.1	3.113537E+07	0.05	3.139198E+07	0.01	2.895182E+07	0.005	1.1177040-00	├ ──-	
Methownshia-		0.15	1.134168E+0/	0.1	1.139067E+07	0.05	1.14118/E+07	0.01	1.09/018E+0/	0.005	0000200	<u> </u>	
Methoyushlas [20]		0.15	1.153144E+07	0.1	1.132936E+07	0.05	1.11782E+07	0.01	1.039432E+07	0.005	9990320	<u> </u>	
Tetrophlers my		0.15	3/10200	0.1	3681211	0.05	3725374	0.01	3691920	0.005	3369360		0.67766670-07
Tatrachiere							<u> </u>	<u> </u>					2.31/330E+0/
i euachioro-m-xyle	աշլույ			1	1	1	1	1		ł		1 0.1	9521665

INITIAL CALIBRATION DATA (Continued)

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03
Calibration:	<u>R9I2904</u>	Instrument:	<u>HP6890-6</u>
		Calibration Date:	09/28/09 00:00

	L	Level 07		Level 08		Level 09		Level 10		Level 11		Level 12	
Compound	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	ng/ul	RF	
Decachlorobiphenyl	0.075	2.285779E+07	0.05	2.278872E+07	0.01	2.423434E+07	0.005	2.574254E+07					
Decachlorobiphenyl [2C]	0.075	8803008	0.05	8978252	0.01	1.018555E+07	0.005	1.101184E+07					
Tetrachloro-m-xylene	0.075	2.569851E+07	0.05	2.500586E+07	0.01	2.328273E+07	0.005	2.38338E+07					
Tetrachloro-m-xylene [2C]	0.075	9526504	0.05	9265848	0.01	8892800	0.005	9276740					

INITIAL CALIBRATION DATA (Continued)

Laboratory:	<u>TestAmerica Buffa</u>	llo		SDC	G:				
Client:	<u>Olin Chlor Alkali I</u>	Products - Cleve	eland, TN	Proj	ect:	<u>Olin - Cha</u>	rles Gibson site	- NY3A9025AE	<u>03759</u>
Calibration:	<u>R912904</u>			Inst	rument:	<u>HP6890-6</u>			
				Cali	bration Date:	<u>09/28/09 0</u>	<u>0:00</u>		
Compound		Mean RF	RF RSD	Mean RT	RT RSD	Linear r ²	Ouad COD	LIMIT	0

Compound	Ivicali IXI	KI KSD	Mican K1	KI KDD	Lincari	Quad COD	Liivii i	× 1
4,4'-DDD	2.19006E+07	9.480439	17.09529	0.0484538	0,99993			
4,4'-DDD [2C]	7962215	6.338822	18.95254	2.454554E-02	0.99998			
4,4'-DDE	2.888943E+07	9.732555	15.77663	4.993852E-02	0.99997			
4,4'-DDE [2C]	1.019583E+07	6.211038	17.67772	3.579122E-02	0.99996			
4,4'-DDT	2.14123E+07	14.36936	17.70203	0.0449234	0.99980			
4,4'-DDT [2C]	7123641	9.815323	19.64778	2.670215E-02	0.99960			
Aldrin	3.082834E+07	8.080935	13.53827	7.117066E-02	0.99998			
Aldrin [2C]	1.166556E+07	6.604578	15.20348	3.524396E-02	0.99998			
alpha-BHC	3.584573E+07	12.73169	10.7546	9.682097E-02	0.99960			
alpha-BHC [2C]	1.37734E+07	12.50096	12.29399	4.348809E-02	0.99991			
alpha-Chlordane	2.961629E+07	4.376397	15.57482	5.533391E-02	0.99995			
alpha-Chlordane [2C]	1.063622E+07	1.669538	17.32776	3.573404E-02	1.00000			
beta-BHC	1.438893E+07	1.759548	11.84956	7.869338E-02	0.99986			
beta-BHC [2C]	5958901	3.657286	13.46889	4.290801E-02	0.99990			
Decachlorobiphenyl	2.362489E+07	5.759798	22.08204	9.490632E-03	0.99995			
Decachlorobiphenyl [2C]	9507167	11.03465	25.21592	6.138726E-03	0.99977			
delta-BHC	3.288152E+07	13.69479	12.2902	7.423003E-02	0.99996			
delta-BHC [2C]	1.286693E+07	11.66999	14.21881	3.917959E-02	0.99992			
Dieldrin	2.965426E+07	8.001368	16.41693	4.713956E-02	0.99996			
Dieldrin [2C]	1.068287E+07	4.79764	18.08091	3.108253E-02	0.99998			
Endosulfan I	2.73306E+07	3.656715	15.87667	5.301915E-02	0.99992			
Endosulfan I [2C]	1.008713E+07	1.662735	17.46332	3.297117E-02	0.99999			
Endosulfan II	2.43636E+07	3.436372	17.42273	5.156096E-02	0.99997			
Endosulfan II [2C]	8941875	1.277227	19.22097	2.289678E-02	1.00000			
Endosulfan sulfate	1.698053E+07	13.59722	19.28673	3.839005E-02	0.99956			
Endosulfan sulfate [2C]	6580221	8.310187	20.56022	1.995262E-02	0.99967			
Endrin	2.533332E+07	9.41182	16.92884	5.368444E-02	0.99999			
Endrin [2C]	8788617	6.254542	18.76311	2.511791E-02	0.99994			
Endrin aldehyde	1.614859E+07	3.742073	18.34277	3.988395E-02	0.99992			
Endrin aldehyde [2C]	6277186	4.77721	19.95123	2.304882E-02	0.99996			
Endrin ketone	2.285868E+07	8.015148	19.89158	3.440551E-02	0.99962			
Endrin ketone [2C]	8534243	2.974258	21.78153	1.607981E-02	0.99970			
gamma-BHC (Lindane)	3.311043E+07	10.40874	11.60176	0.0826149	0.99989			
gamma-BHC (Lindane) [2C]	1.313586E+07	8.881807	13.25124	4.224503E-02	0.99997			
gamma-Chlordane	3.119267E+07	6.321133	15.26992	5.431675E-02	0.99998			
gamma-Chlordane [2C]	1.107746E+07	3.357555	16.9941	3.169023E-02	0.99999			
Heptachlor	3.228413E+07	5.660166	12.80978	7.751327E-02	0.99988			
Heptachlor [2C]	1.277634E+07	4.48319	14.40177	4.018076E-02	0.99999			
Heptachlor epoxide	3.033737E+07	3.475099	14.98502	5.614366E-02	0.99991			
Heptachlor epoxide [2C]	1.129829E+07	1.639064	16.55471	3.298843E-02	0.99999			
Methoxychlor	1.088473E+07	6.063761	18.7498	4.124065E-02	0.99994			
Methoxychlor [2C]	3675613	1.680382	21.08835	1.169048E-02	0.99998			
Tetrachloro-m-xylene	2.471929E+07	4.524952	9.178268	3.654636E-02	0.99991			
Tetrachloro-m-xylene [2C]	9296711	2.783754	10.53044	3.073516E-02	0.99991			

CONTINUING CALIBRATION CHECK

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46176</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	<u>09/25/09</u>
Lab Sample ID:	<u>RI92924-CCV1</u>	Injection Time:	<u>01:59</u>

		CONC. (ng/ul)		RESI	PONSE FACTO	OR	% DIFF / DRIFT		
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)	
4,4'-DDD	L	0.0500	0.0489	2.04568E+07	2.122666E+07		-2.1	15	
4,4'-DDD [2C]	L	0.0500	0.0534	6759941	7514018		6.9	15	
4,4'-DDE	L	0.0500	0.0555	2.359278E+07	2.831326E+07		11.0	15	
4,4'-DDE [2C]	L	0.0500	0.0541	8330355	9450568		8.3	15	
4,4'-DDT	L	0.0500	0.0429	2.103215E+07	1.928204E+07		-14.3	15	
4,4'-DDT [2C]	L	0.0500	0.0451	5960936	5706340		-9.9	15	
Aldrin	L	0.0500	0.0624	2.237609E+07	3.053366E+07		24.7	15 *	
Aldrin [2C]	L	0.0500	0.0547	9926595	1.142506E+07		9.4	15	
alpha-BHC	L	0.0500	0.0607	2.638417E+07	3.61857E+07		21.5	15 *	
alpha-BHC [2C]	L	0.0500	0.0480	1.311296E+07	1.369583E+07		-4.1	15	
alpha-Chlordane	L	0.0500	0.0593	2.322161E+07	2.869424E+07		18.6	15 *	
alpha-Chlordane [2C]	L	0.0500	0.0547	9365786	9911744		9.4	15	
beta-BHC	L	0.0500	0.0610	1.086786E+07	1.357051E+07		22.0	15 *	
beta-BHC [2C]	L	0.0500	0.0516	5499314	5574304		3.3	15	
delta-BHC	L	0.0500	0.0596	2.480519E+07	3.321916E+07		19.2	15 *	
delta-BHC [2C]	L	0.0500	0.0514	1.16449E+07	1.287882E+07		2.9	15	
Dieldrin	L	0.0500	0.0533	2.578433E+07	2.91478E+07		6.7	15	
Dieldrin [2C]	L	0.0500	0.0541	8868618	9911064		8.2	15	
Endosulfan I	L	0.0500	0.0559	2.302769E+07	2.66387E+07		11.9	15	
Endosulfan I [2C]	L	0.0500	0.0546	8448751	9258388		9.2	15	
Endosulfan II	L	0.0500	0.0502	2.255103E+07	2.31825E+07		0.5	15	
Endosulfan II [2C]	L	0.0500	0.0544	7624897	8231508		8.9	15	
Endosulfan sulfate	L	0.0500	0.0463	2.082641E+07	1.958303E+07		-7.4	15	
Endosulfan sulfate [2C]	L	0.0500	0.0517	7253679	7306040		3.3	15	
Endrin	L	0.0500	0.0520	2.399218E+07	2.538876E+07		4.0	15	
Endrin [2C]	L	0.0500	0.0558	7207854	8273568		11.6	15	
Endrin aldehyde	L	0.0500	0.0441	1.714223E+07	1.528342E+07		-11.7	15	
Endrin aldehyde [2C]	L	0.0500	0.0504	5723167	5607404		0.8	15	
Endrin ketone	L	0.0500	0.0457	2.371625E+07	2.23939E+07		-8.6		

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6b46176</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>R192924</u>	Injection Date:	09/25/09
Lab Sample ID:	<u>RI92924-CCV1</u>	Injection Time:	<u>01:59</u>

		CONC	CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)	
Endrin ketone [2C]	L	0.0500	0.0480	8062050	7768144		-3.9		
gamma-BHC (Lindane)	L	0.0500	0.0599	2.463557E+07	3.268134E+07		19.8	15 *	
gamma-BHC (Lindane) [2C]	L	0.0500	0.0497	1.20374E+07	1.272382E+07		-0.7	15	
gamma-Chlordane	L	0.0500	0.0574	2.51191E+07	3.024996E+07		14.8	15	
gamma-Chlordane [2C]	L	0.0500	0.0533	9441841	1.010553E+07		6.5	15	
Heptachlor	L	0.0500	0.0625	2.403694E+07	3.204558E+07		25.0	15 *	
Heptachlor [2C]	L	0.0500	0.0531	1.116489E+07	1.223634E+07		6.2	15	
Heptachlor epoxide	L	0.0500	0.0595	2.378182E+07	2.955908E+07		19.0	15 *	
Heptachlor epoxide [2C]	L	0.0500	0.0553	9485284	1.059097E+07		10.6	15	
Methoxychlor	L	0.0500	0.0409	1.153698E+07	9579822		-18.1	15 *	
Methoxychlor [2C]	L	0.0500	0.0437	3351565	2971964		-12.6	15	

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46177</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	09/25/09
Lab Sample ID:	<u>RI92924-CCV2</u>	Injection Time:	<u>02:35</u>

		CONC. (ng/ul)		RES	RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)	
Decachlorobiphenyl	L	0.0500	0.0477	2.320735E+07	2.164016E+07		-4.6	15	
Decachlorobiphenyl [2C]	L	0.0500	0.0567	7933079	8332668		13,3	15	
Tetrachloro-m-xylene	L	0.0500	0.0697	1.599268E+07	2.349492E+07		39.5	15 *	
Tetrachloro-m-xylene [2C]	L	0.0500	0.0439	9542319	8446886		-12.1	15	

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	<u>TestAmerica Buffalo</u>	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46185</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	09/25/09
Lab Sample ID:	<u>RI92924-CCV3</u>	Injection Time:	<u>07:22</u>

		CONC	. (ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	ТҮРЕ	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
4,4'-DDD	L	0.0500	0.0485	2.04568E+07	2.101408E+07		-3.1	15
4,4'-DDD [2C]	L	0.0500	0.0544	6759941	7648638		8.7	15
4,4'-DDE	L	0.0500	0.0558	2.359278E+07	2.846112E+07		11.5	15
4,4'-DDE [2C]	L	0.0500	0.0551	8330355	9630586		10.3	15
4,4'-DDT	L	0.0500	0.0416	2.103215E+07	1.86918E+07		-16.8	15 *
4,4'-DDT [2C]	L	0.0500	0.0461	5960936	5844112		-7.8	15
Aldrin	L	0.0500	0.0625	2.237609E+07	3.0626E+07		25.1	15 *
Aldrin [2C]	L	0.0500	0.0556	9926595	1.161805E+07		11.1	15
alpha-BHC	L	0.0500	0.0614	2.638417E+07	3.657994E+07		22.7	15 *
alpha-BHC [2C]	L	0.0500	0.0487	1.311296E+07	1.391349E+07		-2.7	15
alpha-Chlordane	L	0.0500	0.0579	2.322161E+07	2.80164E+07		15.8	15 *
alpha-Chlordane [2C]	L	0.0500	0.0556	9365786	1.007173E+07		11.2	15
beta-BHC	L	0.0500	0.0613	1.086786E+07	1.363822E+07		22.6	15 *
beta-BHC [2C]	L	0.0500	0.0523	5499314	5642184		4.5	15
delta-BHC	L	0.0500	0.0596	2.480519E+07	3.324508E+07		19.3	15 *
delta-BHC [2C]	L	0.0500	0.0522	1.16449E+07	1.308444E+07		4.4	15
Dieldrin	L	0.0500	0.0533	2.578433E+07	2.912808E+07		6.6	15
Dieldrin [2C]	L	0.0500	0.0550	8868618	1.008442E+07		10.0	15
Endosulfan I	L	0.0500	0.0560	2,302769E+07	2.665732E+07		12.0	15
Endosulfan I [2C]	L	0.0500	0.0556	8448751	9421694		11.1	15
Endosulfan II	L	0.0500	0.0505	2.255103E+07	2.331326E+07		1.0	15
Endosulfan II [2C]	L	0.0500	0.0553	7624897	8365574		10.6	15
Endosulfan sulfate	L	0.0500	0.0466	2.082641E+07	1.972836E+07		-6.7	15
Endosulfan sulfate [2C]	L	0.0500	0.0528	7253679	7464256		5.6	15
Endrin	L	0.0500	0.0506	2.399218E+07	2.47097E+07		1.2	15
Endrin [2C]	L	0.0500	0.0551	7207854	8174968		10.3	15
Endrin aldehyde	L	0.0500	0.0450	1.714223E+07	1.557378E+07		-10.1	15
Endrin aldehyde [2C]	L	0.0500	0.0516	5723167	5743332		3.3	15
Endrin ketone	L	0.0500	0.0458	2.371625E+07	2.242582E+07		-8.5	

Form Rev: 10/01/2009

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6b46185</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>R192924</u>	Injection Date:	<u>09/25/09</u>
Lab Sample ID:	<u>RI92924-CCV3</u>	Injection Time:	<u>07:22</u>

		CONC. (ng/ul)		RES	RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)	
Endrin ketone [2C]	L	0.0500	0.0495	8062050	8005944		-1.0		
gamma-BHC (Lindane)	L	0.0500	0.0606	2.463557E+07	3.308458E+07		21.2	15 *	
gamma-BHC (Lindane) [2C]	L	0.0500	0.0507	1.20374E+07	1.299561E+07		1.4	15	
gamma-Chlordane	L	0.0500	0.0574	2.51191E+07	3.023404E+07		14.8	15	
gamma-Chlordane [2C]	L	0.0500	0.0542	9441841	1.028133E+07		8.4	15	
Heptachlor	L	0.0500	0.0621	2.403694E+07	3.183798E+07		24.2	15 *	
Heptachlor [2C]	L	0.0500	0.0538	1.116489E+07	1.239748E+07		7.6	15	
Heptachlor epoxide	L	0.0500	0.0597	2.378182E+07	2.96361E+07		19.3	15 *	
Heptachlor epoxide [2C]	L	0.0500	0.0564	9485284	1.079675E+07		12.7	15	
Methoxychlor	L	0.0500	0.0390	1.153698E+07	9131466		-21.9	15 *	
Methoxychlor [2C]	L	0.0500	0.0439	3351565	2985868		-12.2	15	

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46186</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	<u>09/25/09</u>
Lab Sample ID:	<u>RI92924-CCV4</u>	Injection Time:	<u>07:58</u>

		CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Decachlorobiphenyl	L	0.0500	0.0486	2.320735E+07	2.206136E+07		-2.7	15
Decachlorobiphenyl [2C]	L	0.0500	0.0577	7933079	8485864		15.5	15 *
Tetrachloro-m-xylene	L	0.0500	0.0714	1.599268E+07	2.405328E+07		42.7	15 *
Tetrachloro-m-xylene [2C]	L	0.0500	0.0450	9542319	8649184		-10.1	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	<u>TestAmerica Buffalo</u>	SDG:	
Client:	<u> Olin Chlor Alkali Products - Cleveland, TN</u>	Project:	<u>Olin - Charles Gibson site - NY3A9025AE03759</u>
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46196</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	09/25/09
Lab Sample ID:	<u>RI92924-CCV5</u>	Injection Time:	<u>14:07</u>

		CONC.	(ng/ul)	RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
4,4'-DDD	L	0.0500	0.0477	2.04568E+07	2.067614E+07		-4.6	15
4,4'-DDD [2C]	L	0.0500	0.0529	6759941	7437256		5.8	15
4,4'-DDE	L	0.0500	0.0542	2.359278E+07	2.763746E+07		8.5	15
4,4'-DDE [2C]	L	0.0500	0.0475	8330355	8240774		-5.1	15
4,4'-DDT	L	0.0500	0.0422	2.103215E+07	1.896313E+07		-15.6	15 *
4,4'-DDT [2C]	L	0.0500	0.0449	5960936	5680358		-10.3	15
Aldrin	L	0.0500	0.0601	2.237609E+07	2.937326E+07		20.2	15 *
Aldrin [2C]	L	0.0500	0.0535	9926595	1.116237E+07		6.9	15
alpha-BHC	L	0.0500	0.0575	2.638417E+07	3.41269E+07		15.0	15
alpha-BHC [2C]	L	0.0500	0.0470	1.311296E+07	1.341333E+07		-5.9	15
alpha-Chlordane	L	0.0500	0.0539	2.322161E+07	2.600484E+07		7.7	15
alpha-Chlordane [2C]	L	0.0500	0.0485	9365786	8790592		-3.0	15
beta-BHC	L	0.0500	0.0585	1.086786E+07	1.300645E+07		17.1	15 *
beta-BHC [2C]	L	0.0500	0.0505	5499314	5454716		1.0	15
delta-BHC	L	0.0500	0.0574	2.480519E+07	3.193596E+07		14.8	15
delta-BHC [2C]	L	0.0500	0.0504	1.16449E+07	1.261459E+07		0.9	15
Dieldrin	L	0.0500	0.0519	2.578433E+07	2.83377E+07		3.8	15
Dieldrin [2C]	L	0.0500	0.0539	8868618	9867552		7.7	15
Endosulfan I	L	0.0500	0.0542	2.302769E+07	2.579374E+07		8.4	15
Endosulfan I [2C]	L	0.0500	0.0542	8448751	9192494		8.5	15
Endosulfan II	L	0.0500	0.0492	2.255103E+07	2.26884E+07		-1,6	15
Endosulfan II [2C]	L	0.0500	0.0536	7624897	8103122		7.2	15
Endosulfan sulfate	L	0.0500	0.0453	2.082641E+07	1.913686E+07		-9.5	15
Endosulfan sulfate [2C]	L	0.0500	0.0510	7253679	7218568		2.1	15
Endrin	L	0.0500	0.0493	2.399218E+07	2.403744E+07		-1.5	15
Endrin [2C]	L	0.0500	0.0538	7207854	7975658		7.7	15
Endrin aldehyde	L	0.0500	0.0434	1.714223E+07	1.502319E+07		-13.2	15
Endrin aldehyde [2C]	L	0.0500	0.0488	5723167	5432572		-2.4	15
Endrin ketone	L	0.0500	0.0450	2.371625E+07	2.202136E+07		-10.1	<u> </u>

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CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6b46196</u>	Calibration Date:	07/23/09 10:05
Sequence:	<u>RI92924</u>	Injection Date:	<u>09/25/09</u>
Lab Sample ID:	<u>RI92924-CCV5</u>	Injection Time:	14:07

		CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Endrin ketone [2C]	L	0.0500	0.0482	8062050	7799816		-3.5	
gamma-BHC (Lindane)	L	0.0500	0.0576	2.463557E+07	3.135132E+07		15.2	15 *
gamma-BHC (Lindane) [2C]	L	0.0500	0.0489	1.20374E+07	1.250498E+07		-2.3	15
gamma-Chlordane	L	0.0500	0.0557	2.51191E+07	2.93017E+07		11.3	15
gamma-Chlordane [2C]	L	0.0500	0.0544	9441841	1.032827E+07		8.8	15
Heptachlor	L	0.0500	0.0605	2.403694E+07	3.09852E+07		21.0	15 *
Heptachlor [2C]	L	0.0500	0.0526	1.116489E+07	1.211289E+07		5.2	15
Heptachlor epoxide	L	0.0500	0.0577	2.378182E+07	2.863448E+07		15.4	15 *
Heptachlor epoxide [2C]	L	0.0500	0.0490	9485284	9364814		-2.0	15
Methoxychlor	L	0.0500	0.0399	1.153698E+07	9327202		-20.3	15 *
Methoxychlor [2C]	L	0.0500	0.0458	3351565	3113628		-8.4	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9H1803</u>
Lab File ID:	<u>6A46197</u>	Calibration Date:	<u>07/23/09 10:05</u>
Sequence:	<u>R192924</u>	Injection Date:	09/25/09
Lab Sample ID:	<u>R192924-CCV6</u>	Injection Time:	<u>14:53</u>

		CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Decachlorobiphenyl	L	0.0500	0.0490	2.320735E+07	2.224194E+07		-1.9	15
Decachlorobiphenyl [2C]	L	0.0500	0.0579	7933079	8506102		15.8	15 *
Tetrachloro-m-xylene	L	0.0500	0.0673	1.599268E+07	2.263902E+07		34.5	15 *
Tetrachloro-m-xylene [2C]	L	0.0500	0.0453	9542319	8721640		-9.3	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

Laboratory:	<u>TestAmerica Buffalo</u>	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
Lab File ID:	<u>6A46228</u>	Calibration Date:	09/28/09_00:00
Sequence:	<u>RI92925</u>	Injection Date:	09/28/09
Lab Sample ID:	<u>RI92925-CCV1</u>	Injection Time:	<u>19:21</u>

		CONC	. (ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
4,4'-DDD	L	0.0500	0.0510	2.19006E+07	2.365576E+07		2.1	15
4,4'-DDD [2C]	L	0.0500	0.0514	7962215	8510182		2.8	15
4,4'-DDE	L	0.0500	0.0504	2.888943E+07	3.085476E+07		0.8	15
4,4'-DDE [2C]	L	0.0500	0.0511	1.019583E+07	1.084E+07		2.3	15
4,4'-DDT	L	0.0500	0.0490	2.14123E+07	2.284726E+07		-2.1	15
4,4'-DDT [2C]	L	0.0500	0.0507	7123641	7680412		1.4	15
Aldrin	L	0.0500	0.0511	3.082834E+07	3.302702E+07		2.1	15
Aldrin [2C]	L	0.0500	0.0512	1.166556E+07	1.243099E+07		2.3	15
alpha-BHC	L	0.0500	0.0499	3.584573E+07	3.850648E+07		-0.2	15
alpha-BHC [2C]	L	0.0500	0.0500	1.37734E+07	1.484582E+07		-0.09	15
alpha-Chlordane	L	0.0500	0.0506	2.961629E+07	3.076536E+07		1.3	15
alpha-Chlordane [2C]	L	0.0500	0.0519	1.063622E+07	1.109377E+07		3.9	15
beta-BHC	L	0.0500	0.0508	1.438893E+07	1.454043E+07		1.5	15
beta-BHC [2C]	L	0.0500	0.0520	5958901	6058968		4.0	15
delta-BHC	L	0.0500	0.0499	3.288152E+07	3.560252E+07		-0.2	15
delta-BHC [2C]	L	0.0500	0.0505	1.286693E+07	1.39696E+07		1.1	15
Dieldrin	L	0.0500	0.0507	2.965426E+07	3.15207E+07		1.3	15
Dieldrin [2C]	L	0.0500	0.0512	1.068287E+07	1.126862E+07		2.4	15
Endosulfan I	L	0.0500	0.0512	2.73306E+07	2.851996E+07		2.3	15
Endosulfan I [2C]	L	0.0500	0.0520	1.008713E+07	1.054195E+07		4.0	15
Endosulfan II	L	0.0500	0.0508	2.43636E+07	2.526432E+07		1.6	15
Endosulfan II [2C]	L	0.0500	0.0518	8941875	9261038		3.6	15
Endosulfan sulfate	L	0.0500	0.0501	1.698053E+07	1.848313E+07		0.2	15
Endosulfan sulfate [2C]	L	0.0500	0.0518	6580221	7186258		3.7	15
Endrin	L	0.0500	0.0500	2.533332E+07	2.679058E+07		-0.06	15
Endrin [2C]	L	0.0500	0.0504	8788617	9218864	,	0.9	15
Endrin aldehyde	L	0.0500	0.0510	1.614859E+07	1.681239E+07		2.0	15
Endrin aldehyde [2C]	L	0.0500	0.0521	6277186	6354612		4.3	15
Endrin ketone	L	0.0500	0.0503	2.285868E+07	2.417048E+07		0.6	

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R912904</u>
Lab File ID:	<u>6b46228</u>	Calibration Date:	09/28/09 00:00
Sequence:	<u>R192925</u>	Injection Date:	09/28/09
Lab Sample ID:	<u>RI92925-CCV1</u>	Injection Time:	<u>19:21</u>

		CONC.	(ng/ul)	RESPONSE FACTOR		% DIFF / DRIFT		
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Endrin ketone [2C]	L	0.0500	0.0524	8534243	9114926		4.9	
gamma-BHC (Lindane)	L	0.0500	0.0499	3.311043E+07	3.514842E+07		-0.2	15
gamma-BHC (Lindane) [2C]	L	0.0500	0.0506	1.313586E+07	1.402227E+07		1.1	15
gamma-Chlordane	L	0.0500	0.0506	3.119267E+07	3.278444E+07		1.2	15
gamma-Chlordane [2C]	L	0.0500	0.0518	1.107746E+07	1.169853E+07		3.6	15
Heptachlor	L	0.0500	0.0513	3.228413E+07	3.421402E+07		2.7	15
Heptachlor [2C]	L	0.0500	0.0510	1.277634E+07	1.339778E+07		2.1	15
Heptachlor epoxide	L	0.0500	0.0516	3.033737E+07	3.185516E+07		3.2	15
Heptachlor epoxide [2C]	L	0.0500	0.0521	1.129829E+07	1.182572E+07		4.1	15
Methoxychlor	L	0.0500	0.0497	1.088473E+07	1.123666E+07		-0.5	15
Methoxychlor [2C]	L	0.0500	0.0534	3675613	3953120		6.9	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	<u>Olin Chlor Alkali Products - Cleveland, TN</u>	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R912904</u>
Lab File ID:	<u>6A46229</u>	Calibration Date:	09/28/09 00:00
Sequence:	<u>RI92925</u>	Injection Date:	<u>09/28/09</u>
Lab Sample ID:	<u>RI92925-CCV2</u>	Injection Time:	<u>19:57</u>

		CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Decachlorobiphenyl	L	0.0500	0.0501	2.362489E+07	2.28194E+07		0.1	15
Decachlorobiphenyl [2C]	L	0.0500	0.0510	9507167	9000380		2.1	15
Tetrachloro-m-xylene	L	0.0500	0.0486	2.471929E+07	2.476018E+07		-2.7	15
Tetrachloro-m-xylene [2C]	L	0.0500	0.0484	9296711	9133898		-3.2	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
Lab File ID:	<u>6A46238</u>	Calibration Date:	09/28/09 00:00
Sequence:	<u>R192925</u>	Injection Date:	09/29/09
Lab Sample ID:	<u>RI92925-CCV3</u>	Injection Time:	<u>01:20</u>

		CONC.	(ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
4,4'-DDD	L	0.0500	0.0500	2.19006E+07	2.314924E+07		-0.06	15
4,4'-DDD [2C]	L	0.0500	0.0496	7962215	8210372		-0.8	15
4,4'-DDE	L	0.0500	0.0488	2.888943E+07	2.98816E+07		-2.3	15
4,4'-DDE [2C]	L	0.0500	0.0488	1.019583E+07	1.032138E+07		-2.5	15
4,4'-DDT	L	0.0500	0.0467	2.14123E+07	2.171988E+07		-6.6	15
4,4'-DDT [2C]	L	0.0500	0.0481	7123641	7267228		-3.8	15
Aldrin	L	0.0500	0.0508	3.082834E+07	3.287906E+07		1.7	15
Aldrin [2C]	L	0.0500	0.0497	1.166556E+07	1.207813E+07		-0.5	15
alpha-BHC	L	0.0500	0.0495	3.584573E+07	3.816192E+07		-1.1	15
alpha-BHC [2C]	L	0.0500	0.0496	1.37734E+07	1.47447E+07		-0.7	15
alpha-Chlordane	L	0.0500	0.0485	2.961629E+07	2.94643E+07		-3.0	15
alpha-Chlordane [2C]	L	0.0500	0.0494	1.063622E+07	1.055463E+07		-1.2	15
beta-BHC	L	0.0500	0.0505	1.438893E+07	1.446126E+07		0.9	15
beta-BHC [2C]	L	0.0500	0.0515	5958901	5998370		3.0	15
delta-BHC	L	0.0500	0.0504	3.288152E+07	3.59615E+07		0.8	15
delta-BHC [2C]	L	0.0500	0.0503	1.286693E+07	1.391531E+07		0.7	15
Dieldrin	L	0.0500	0.0487	2.965426E+07	3.02615E+07		-2.6	15
Dieldrin [2C]	L	0.0500	0.0487	1.068287E+07	1.071194E+07		-2.5	15
Endosulfan I	L	0.0500	0.0498	2.73306E+07	2.775662E+07		-0.4	15
Endosulfan I [2C]	L	0.0500	0.0497	1.008713E+07	1.00671E+07		-0.7	15
Endosulfan II	L	0.0500	0.0486	2.43636E+07	2.41474E+07		-2.8	15
Endosulfan II [2C]	L	0.0500	0.0493	8941875	8811060		-1.4	15
Endosulfan sulfate	L	0.0500	0.0527	1.698053E+07	1.949542E+07		5.4	15
Endosulfan sulfate [2C]	L	0.0500	0.0541	6580221	7516648		8.3	15
Endrin	L	0.0500	0.0473	2.533332E+07	2.534406E+07		-5.3	15
Endrin [2C]	L	0.0500	0.0463	8788617	8444144		-7.4	15
Endrin aldehyde	L	0.0500	0.0485	1.614859E+07	1.599357E+07		-2.9	15
Endrin aldehyde [2C]	L	0.0500	0.0490	6277186	5976226		-2.0	15
Endrin ketone	L	0.0500	0.0505	2.285868E+07	2.425088E+07		0.9	

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
Lab File ID:	<u>6b46238</u>	Calibration Date:	09/28/09 00:00
Sequence:	<u>R192925</u>	Injection Date:	09/29/09
Lab Sample ID:	<u>RI92925-CCV3</u>	Injection Time:	<u>01:20</u>

	1 1	CONC	. (ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Endrin ketone [2C]	L	0.0500	0.0523	8534243	9090472		4.6	
gamma-BHC (Lindane)	L	0.0500	0.0498	3.311043E+07	3.509036E+07		-0.3	15
gamma-BHC (Lindane) [2C]	L	0.0500	0.0503	1.313586E+07	1.393859E+07		0.5	15
gamma-Chlordane	L	0.0500	0.0489	3.119267E+07	3.16869E+07		-2.2	15
gamma-Chlordane [2C]	L	0.0500	0.0492	1.107746E+07	1.109796E+07		-1.7	15
Heptachlor	L	0.0500	0.0525	3.228413E+07	3.498288E+07		5.0	15
Heptachlor [2C]	L	0.0500	0.0512	1.277634E+07	1.342811E+07		2.3	15
Heptachlor epoxide	L	0.0500	0.0503	3.033737E+07	3.103558E+07		0.5	15
Heptachlor epoxide [2C]	L	0.0500	0.0501	1.129829E+07	1.137175E+07		0.1	15
Methoxychlor	L	0.0500	0.0478	1.088473E+07	1.077538E+07		-4.5	15
Methoxychlor [2C]	L	0.0500	0.0500	3675613	3702524		0.09	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

TestAmerica Buffalo	SDG:	
Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
<u>6A46239</u>	Calibration Date:	09/28/09_00:00
<u>RI92925</u>	Injection Date:	09/29/09
<u>RI92925-CCV4</u>	Injection Time:	<u>01:56</u>
	TestAmerica Buffalo Olin Chlor Alkali Products - Cleveland, TN HP6890-6 6A46239 R192925 R192925-CCV4	TestAmerica BuffaloSDG:Olin Chlor Alkali Products - Cleveland, TNProject:HP6890-6Calibration:6A46239Calibration Date:R192925Injection Date:R192925-CCV4Injection Time:

		CONC	C. (ng/ul)	RES	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Decachlorobiphenyl	L	0.0500	0.0486	2.362489E+07	2.216634E+07		-2.8	15
Decachlorobiphenyl [2C]	L	0.0500	0.0484	9507167	8548204		-3.3	15
Tetrachloro-m-xylene	L	0.0500	0.0476	2.471929E+07	2.423898E+07		-4.7	15
Tetrachloro-m-xylene [2C]	L	0.0500	0.0496	9296711	9366856		-0.8	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

Laboratory:	<u>TestAmerica Buffalo</u>	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	<u>Olin - Charles Gibson site - NY3A9025AE03759</u>
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
Lab File ID:	<u>6A46246</u>	Calibration Date:	<u>09/28/09 00:00</u>
Sequence:	<u>RI92925</u>	Injection Date:	<u>09/29/09</u>
Lab Sample ID:	<u>R192925-CCV5</u>	Injection Time:	<u>06:08</u>
Lab File ID: Sequence: Lab Sample ID:	<u>HP6890-6</u> <u>6A46246</u> <u>R192925</u> <u>R192925-CCV5</u>	Calibration: Calibration Date: Injection Date: Injection Time:	<u>K912904</u> 09/28/09 00:00 09/29/09 06:08

		CONC	. (ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
4,4'-DDD	L	0.0500	0.0549	2.19006E+07	2.550408E+07		9.9	15
4,4'-DDD [2C]	L	0.0500	0.0553	7962215	9177292		10.6	15
4,4'-DDE	L	0.0500	0.0512	2.888943E+07	3.134364E+07		2.3	15
4,4'-DDE [2C]	L	0.0500	0.0515	1.019583E+07	1.092511E+07		3.1	15
4,4'-DDT	L	0.0500	0.0432	2.14123E+07	2.0024E+07		-13.5	15
4,4'-DDT [2C]	L	0.0500	0.0451	7123641	6798776		-9.7	15
Aldrin	L	0.0500	0.0533	3.082834E+07	3.453108E+07		6.7	15
Aldrin [2C]	L	0.0500	0.0524	1.166556E+07	1.274089E+07		4.8	15
alpha-BHC	L	0.0500	0.0506	3.584573E+07	3.904752E+07		1.2	15
alpha-BHC [2C]	L	0.0500	0.0525	1.37734E+07	1.563132E+07		5.0	15
alpha-Chlordane	L	0.0500	0.0520	2.961629E+07	3.15811E+07		4.0	15
alpha-Chlordane [2C]	L	0.0500	0.0524	1.063622E+07	1.118817E+07		4.7	15
beta-BHC	L	0.0500	0.0522	1.438893E+07	1.494139E+07		4.3	15
beta-BHC [2C]	L	0.0500	0.0541	5958901	6296032		8.2	15
delta-BHC	L	0.0500	0.0523	3.288152E+07	3.739344E+07]	4.6	15
delta-BHC [2C]	L	0.0500	0.0530	1.286693E+07	1.466839E+07		5.9	15
Dieldrin	L	0.0500	0.0513	2.965426E+07	3.19026E+07		2.5	15
Dieldrin [2C]	L	0.0500	0.0516	1.068287E+07	1.136104E+07		3.3	15
Endosulfan I	L	0.0500	0.0521	2.73306E+07	2.903426E+07		4.2	15
Endosulfan I [2C]	L	0.0500	0.0525	1.008713E+07	1.063605E+07		4.9	15
Endosulfan II	L	0.0500	0.0511	2.43636E+07	2.54027E+07		2.2	15
Endosulfan II [2C]	L	0.0500	0.0522	8941875	9324784		4.3	15
Endosulfan sulfate	L	0.0500	0.0558	1.698053E+07	2.072356E+07		11.7	15
Endosulfan sulfate [2C]	L	0.0500	0.0578	6580221	8038458	1	15.5	15 *
Endrin	L	0.0500	0.0501	2.533332E+07	2.683624E+07		0.1	15
Endrin [2C]	L	0.0500	0.0504	8788617	9205388		0.7	15
Endrin aldehyde	L	0.0500	0.0507	1.614859E+07	1.672368E+07		1.5	15
Endrin aldehyde [2C]	L	0.0500	0.0519	6277186	6320974		3.7	15
Endrin ketone	L	0.0500	0.0527	2.285868E+07	2.535352E+07		5.3	

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R9I2904</u>
Lab File ID:	<u>6b46246</u>	Calibration Date:	09/28/09 00:00
Sequence:	<u>R192925</u>	Injection Date:	09/29/09
Lab Sample ID:	<u>RI92925-CCV5</u>	Injection Time:	<u>06:08</u>

		CONC.	(ng/ul)	RESI	PONSE FACTO	OR	% DIFF	/ DRIFT
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Endrin ketone [2C]	L	0.0500	0.0551	8534243	9595078		10.3	
gamma-BHC (Lindane)	L	0.0500	0.0513	3.311043E+07	3.612664E+07		2.5	15
gamma-BHC (Lindane) [2C]	L	0.0500	0.0528	1.313586E+07	1.465889E+07		5.6	15
gamma-Chlordane	L	0.0500	0.0516	3.119267E+07	3.343436E+07		3.1	15
gamma-Chlordane [2C]	L	0.0500	0.0519	1.107746E+07	1.172186E+07		3.8	15
Heptachlor	L	0.0500	0.0541	3.228413E+07	3.609508E+07		8.3	15
Heptachlor [2C]	L	0.0500	0.0531	1.277634E+07	1.395266E+07		6.3	15
Heptachlor epoxide	L	0.0500	0.0528	3.033737E+07	3.261676E+07		5.7	15
Heptachlor epoxide [2C]	L	0.0500	0.0526	1.129829E+07	1.195736E+07		5.3	15
Methoxychlor	L	0.0500	0.0465	1.088473E+07	1.049475E+07		-6.9	15
Methoxychlor [2C]	L	0.0500	0.0497	3675613	3674044		-0.7	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

CONTINUING CALIBRATION CHECK

8081A

Laboratory:	TestAmerica Buffalo	SDG:	
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Instrument ID:	<u>HP6890-6</u>	Calibration:	<u>R912904</u>
Lab File ID:	<u>6A46247</u>	Calibration Date:	<u>09/28/09 00:00</u>
Sequence:	<u>RI92925</u>	Injection Date:	<u>09/29/09</u>
Lab Sample ID:	<u>R192925-CCV6</u>	Injection Time:	<u>06:44</u>

		CONC. (ng/ul)		RESPONSE FACTOR			% DIFF / DRIFT	
COMPOUND	TYPE	STD	CCV	ICAL	CCV	MIN (#)	CCV	LIMIT (#)
Decachlorobiphenyl	L	0.0500	0.0512	2.362489E+07	2.332846E+07		2.4	15
Decachlorobiphenyl [2C]	L	0.0500	0.0514	9507167	9055184		2.7	15
Tetrachloro-m-xylene	L	0.0500	0.0481	2.471929E+07	2.44811E+07		-3.8	15
Tetrachloro-m-xylene [2C]	L	0.0500	0.0517	9296711	9769592		3.4	15

Column to be used to flag Response Factor and %Diff/Drift values with an asterisk

* Values outside of QC limits

Calibration Type Legend:

A: Average RF

L: Linear through Intercept

Q: Quadratic

L0: Linear forced through Zero

L1: 1/x Weighted Linear through Intercept

L2: 1/x2 Weighted Linear through Intercept

L01: 1/x Weighted Linear forced through Zero

ANALYSIS BATCH (SEQUENCE) SUMMARY

Laboratory:	TestAmerica B	uffalo	SDG:			
Client: Olin Chlor Alkali Products		ali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE0375		
Sequence:	<u>R192924</u>		Instrument:	<u>HP6890-6</u>		
			Calibration:	<u>R9H1803</u>		
Sample Name		Lab Sample ID	Lab File ID	Analysis Date/Time		
Calibration Chec	k	RI92924-CCV1	6A46176	09/25/09 01:59		
Calibration Chec	k	RI92924-CCV1	6b46176	09/25/09 01:59		
Calibration Chec	k	RI92924-CCV2	6A46177	09/25/09 02:35		
Calibration Chec	k	RI92924-CCV2	6b46177	09/25/09 02:35		
LCS		9120007-BS1	6A46181	09/25/09 04:58		
LCS		9120007-BS1	6b46181	09/25/09 04:58		
Calibration Chec	k	RI92924-CCV3	6A46185	09/25/09 07:22		
Calibration Chec	k	R192924-CCV3	6b46185	09/25/09 07:22		
Calibration Chec	k	RI92924-CCV4	6A46186	09/25/09 07:58		
Calibration Chec	k	RI92924-CCV4	6b46186	09/25/09 07:58		
Blank		9I20007-BLK1	6A46193	09/25/09 12:09		
Blank		9I20007-BLK1	6b46193	09/25/09 12:09		
Calibration Chec	k	RI92924-CCV5	6A46196	09/25/09 14:07		
Calibration Chec	k	RI92924-CCV5	6b46196	09/25/09 14:07		
Calibration Chec	k	RI92924-CCV6	6A46197	09/25/09 14:53		
Calibration Chec	k	RI92924-CCV6	6b46197	09/25/09 14:53		

ANALYSIS BATCH (SEQUENCE) SUMMARY

Laboratory:	TestAmerica B	<u>uffalo</u>	SDG:	
Client:	Olin Chlor Alk	ali Products - Cleveland, TN	Project:	Olin - Charles Gibson site - NY3A9025AE03759
Sequence:	<u>RI92925</u>		Instrument:	<u>HP6890-6</u>
			Calibration:	<u>R912904</u>
Sample Name		Lab Sample ID	Lab File ID	Analysis Date/Time
Performance Mi	ix	RI92925-PEM1	6A46227	09/28/09 18:44
Performance Mi	ix	RI92925-PEM1	6b46227	09/28/09 18:44
Calibration Che	ck	RI92925-CCV1	6A46228	09/28/09 19:21
Calibration Che	ck	RI92925-CCV1	6b46228	09/28/09 19:21
Calibration Che	ck	RI92925-CCV2	6A46229	09/28/09 19:57
Calibration Che	ck	RI92925-CCV2	6b46229	09/28/09 19:57
Calibration Che	ck	RI92925-CCV3	6A46238	09/29/09 01:20
Calibration Che	ck	RI92925-CCV3	6b46238	09/29/09 01:20
Calibration Che	ck	RI92925-CCV4	6A46239	09/29/09 01:56
Calibration Che	ck	RI92925-CCV4	6b46239	09/29/09 01:56
CGS-US1-0917	709	RSI0634-01	6A46240	09/29/09 02:32
CGS-US1-0917	709	RS10634-01	6b46240	09/29/09 02:32
CGS-USD1-09	1709	RSI0634-02	6A46241	09/29/09 03:08
CGS-USD1-09	1709	RSI0634-02	6b46241	09/29/09 03:08
CGS-DS1-0917	709	RSI0634-03	6A46242	09/29/09 03:44
CGS-DS1-0917	709	RSI0634-03	6b46242	09/29/09 03:44
CGS-DSD1-09	1709	RSI0634-04	6A46243	09/29/09 04:20
CGS-DSD1-09	1709	RSI0634-04	6b46243	09/29/09 04:20
Calibration Che	ck	RI92925-CCV5	6A46246	09/29/09 06:08
Calibration Che	ck	RI92925-CCV5	6b46246	09/29/09 06:08
Calibration Che	ck	RI92925-CCV6	6A46247	09/29/09 06:44
Calibration Che	ck	RI92925-CCV6	6b46247	09/29/09 06:44

Form 10A IDENTIFICATION SUMMARY FOR SINGLE COMPONENT ANALYTES

DRAFT: CGS-US1-091709

Lab Name: TestAmerica Buffalo					SDG No.:	DRAFT		
Client: Olin Chlor Alkali Products - Cleveland, TN				٧٧	Project: Olin - Charles Gibson site - NY3A9025AE03759			9025AE03759
Lab Sample ID: RSI0634-01				Date(s) Analyzed:		09/29/2009	09/29/2009	
Instrument ID (1): HP6890-6Column 1		_	Instrument ID (HP6890-6Column 2			
GC Column (1):		D:	(mm) GC Column (2)		ID:		(mm)	
AN	LYTE	COL	RT	RT WINDOW		CONCENTRATION	%D]
				FROM	то			
alpha-BHC		1	10.76	10.70	10.80	240		
		2	12.30	12.24	12.34	270	13	
		2	13.48	13.42	13.52	280	7	
beta-BHC		1	11.86	11.80	11.90	260		
delta-BHC		1	12.30	12.24	12.34	39		
		2	14.23	14.17	14.27	42	8	
gamma-BHC (Lindane)		1	11.61	11.55	11.65	18		
		2	13.26	13.20	13.30	26	41	
PRAFT: CGS-USD1-091709

Lab Name:	TestAmerica Buffal	0			SDG No.:			
Client:	Olin Chlor Alkali Pr	roducts - C	Cleveland, TN	٧٧	Project:	site - NY3A	9025AE03759	
Lab Sample ID:	RSI0634-02			Date	Date(s) Analyzed: 09/29/2009 0			
Instrument ID (1):	HP6890-6Column 1		-	Ir	istrument ID (HP6890-6Column 2		
GC Column (1):	Ι	D:	(mm)	G	C Column (2):	IE):	(mm)
AN	ALYTE	COL	RT	RT W	INDOW	CONCENTRATION	%D]
				FROM	то]
alpha-BHC		1	10.77	10.70	10.80	530		
		2	12.31	12.24	12.34	560	7	
		2	13.48	13.42	13.52	200	4	
beta-BHC		1	11.86	11.80	11.90	190		
delta-BHC		1	12.30	12.24	12.34	36		
		2	14.23	14.17	14.27	39	7	
gamma-BHC (I	Lindane)	1	11.61	11.55	11.65	19		
		2	13.26	13.20	13.30	28	46	

DRAFT: CGS-DS1-091709

	Lab Name:	TestAmerica Buffal	0			SDG No.:	DRAFT		
	Client:	Olin Chlor Alkali Pi	Din Chlor Alkali Products - Cleveland, TN				Project: Olin - Charles Gibson site - NY3A9025AE037		
	Lab Sample ID:	RS10634-03			Date	e(s) Analyzed:	09/29/2009	09/29/2009	
In	strument ID (1):	HP6890-6Column 1		_	In	istrument ID (HP6890-6Column 2		
	GC Column (1):	I	D:	(mm)	GC	C Column (2):	Ι	D:	(mm)
	ANA	ALYTE	COL	RT	RT WI	NDOW	CONCENTRATION	%D] .
					FROM	TO			· ·
	alpha-BHC		1	10.76	10.70	10.80	210		
			2	12.30	12.24	12.34	240	12	
			2	13.48	13.42	13.52	69	5	
	beta-BHC		1	11.86	11.80	11.90	73]
	delta-BHC		1	12.30	12.24	12.34	32]
			2	14.22	14.17	14.27	34	8]

PRAFT: CGS-DSD1-091709

Lab Name:	TestAmerica Buffal	0			SDG No.:	DRAFT		
Client:	Olin Chlor Alkali Pr	Cleveland, Th	NN	Project	Olin - Charles Gibson	Olin - Charles Gibson site - NY3A9025AE03759		
Lab Sample ID:	RSI0634-04			Date	e(s) Analyzed:	09/29/2009 0	9/29/2009	
instrument ID (1):	HP6890-6Column 1		_	Ir	strument ID (HP6890-6Column 2		
GC Column (1):	I	D:	(mm)	G	C Column (2):	ID	:	(mm)
AN	ALYTE	COL	RT	RT WI	INDOW	CONCENTRATION	%D]
				FROM	TO			
alpha-BHC		1	10.76	10.70	10.80	93		
		2	12.30	12.24	12.34	110	18]
		2	13.48	13.42	13.52	180	7]
beta-BHC		1	11.86	11.80	11.90	170]
delta-BHC		1	12.30	12.24	12.34	24		
		2	14.23	14.17	14.27	26	8	

DRAFT: LCS

	Lab Name:	TestAmerica Buffal	0			SDG No.:	DRAFT			
	Client:	Olin Chlor Alkali Pi	Cleveland, TN	1	Project	Olin - Charles Gibson	site - NY3A	9025AE03759		
	Lab Sample ID:	9I20007-BS1			Date	(s) Analyzed:	09/25/2009 0	9/25/2009		
In	strument ID (1):	HP6890-6Column 1		_	In	strument ID (HP6890-6Column 2			
	GC Column (1):	I	D:	(mm)	GC	Column (2):	ID	:	(mm)	
	AN	ALYTE	COL	RT	RT WI	NDOW	CONCENTRATION	%D		
	alvha-BHC			10.94	FROM	10 10,99 11,27	14.9			
			2	12.48	12.7 6	12.86	13.1	14		
			2	13.65	13. 94	13.71 1 4.04 -	14.6	14		
	beta-BHC		1	12.03	12.27 %	12.378	16.7			
	delta-BHC		1	12.48	12. 72 43	12.8 2 53	16.0			
			2	14.41	14.7 0	14.80	14.8	8		
	gamma-BHC (I	Lindane)	1	11.79	12.03	12,13,84	15.6			
			2	13.44	13.739	13.83-9	13.8	13		

AU 10/1/09

TotalChrom Method File H:\TU	RI	306\6890)-06\	6a-(07-23-09).mth
Printed by	:	BescoD	on:	07/24/2009	07:36:15
Created by	;	BescoD	on:	07/24/2009	06:58:40
Edited by	;	BescoD	on:	07/24/2009	07:36:10
Number of Times Edited	:	5			
Number of Times Calibrated	:	2651			
Description: CURVE 07-23-0)9				

Global Sample Information

Default Sample Volume	:	1.000 ul
Quantitation Units	;	ng
Void Time	:	0.000 min
Correct amounts during calibration	:	Yes
Convert unknowns to concentration units	:	Yes
Reject outliers during calibration	:	No

An External Standard calibration will be used Unknown peaks will be quantitated using a response factor of 1.000000e+06 First peak will be relative retention reference

Component Information

alpha-BHC Component Type : Single Peak Component Retention Time : 11.233 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order

Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates

А	0.0050	104357.20	32529.05		 1
В	0.0100	217038.10	67803.66		 1
С	0.0500	1422693.20	454715.04	************	 1
D	0.1000	2951163.20	927945.19	**************	 1
Ε	0.1500	4707018.20	1.46e+06		 1

Calibration Curve : $y = (-109541.526833) + (31587229.114763)x + (0.000000)x^{2} + (0.000000)x^{3}$ R-squared : 0.998609

gamma-BHC

Component Type: Single Peak ComponentRetention Time: 12.085 minSearch Window: 5.00 s, 0.00 %Reference Component :Find peak closest to expected RT in windowCalibrating Area versus Amount using a 1st Order FitCurve will ignore the originAmounts will not be scaled prior to the regressionWeighting factor for the regression: 1Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Processed by: <u>DYB</u> 7 <u>B4107</u> Reviewed by: <u>DYY</u> 8 124105

Cal Lev	libration L vel Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	;
A B C D E		0.0050 0.0100 0.0500 0.1000 0.1500	102120.20 208946.50 1313185.60 2698991.50 4290828.70	30487.42 62929.01 403167.56 827260.29 1.31e+06			1 1 1 1 1 1	
Cal R-s	libration C squared	Curve:y	= (-87595.37 .998757	6760) + (28	736663.704	523)x + (0.00	0000)x^2 + ((0.000000)x^3
beta- Coi Rei Sea Fin Cal Cui Am We Coi	BHC mponent i tention Tii arch Winc ference C d peak clu librating A rve will ign ounts will sighting fa mponent	Type me low omponen osest to e vrea versu nore the c l not be so ctor for th standard	: Single Pe : 12.330 m : 5.00 s, 0. t : xpected RT in s Amount us origin caled prior to re regression: purity percent	eak Compon in .00 % ing a 1st Or the regressi : 1 tage : 100.0	lent der Fit on 000%			
Use L V V V V	er Values .abel : /alue 1 : /alue 2 : /alue 3 : /alue 4 : /alue 5 :	0.020000 0.000000 0.000000 0.000000 0.000000)))					
Cal Lev	libration L vel Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	3
A B C D E		0.0050 0.0100 0.0500 0.1000 0.1500	53416.10 103633.40 551638.90 1086215.80 1709670.80	15497.55 30429.29 164775.79 325971.18 513910.88			1	
Cal R-s	libration C squared	Curve:y	= (-13068.51 .999090	8682) + (11	333071.324	368)x + (0.00	0000)x^2 + (i	0.000000)x^3
delta Coi Rei Sea Fin Cal Cui Am We Coi	R-squared : 0.999090 delta-BHC Component Type : Single Peak Component Retention Time : 12.776 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%							
Us L V V V V	er Values .abel : /alue 1 : /alue 2 : /alue 3 : /alue 4 : /alue 5 :	0.020000 0.000000 0.000000 0.000000 0.000000)))					
Cal Lev	libration L vel Name	.evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	;
A B C D E		0.0050 0.0100 0.0500 0.1000 0.1500	99325.50 205191.00 1331457.80 2754822.00 4419651.26	29585.29 61484.26 404736.00 838079.11 1.33e+06			1	
Cal R-s	libration C squared	Curve:y :0	= (-103556.2 .998328	239555) + (2	9613424.220	0384)x + (0.0	00000)x^2 +	(0.000000)x^3

Heptachlor Component Type : Single Peak Component Retention Time : 13.299 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area Height 107993.00 31928.31 ----- ---1 0.0050 А 0.0100 215243.70 64186.05 -----в 1 С 0.0500 1267228.80 386320.86 -----1 D 1 1 E 0.1500 3948257.70 1.18e+06 -------Calibration Curve : y = (-47233.435424) + (26395032.019260)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999562 Aldrin : Single Peak Component Component Type Retention Time : 14.034 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area Height А 0.0050 95257.00 28243.18 -----1 0.0100 193158.60 57786.96 -----1 В 0.0500 1194589.00 363472.70 ------1 С D 0.1000 2421194.40 727111.12 ------1 0.1500 3811422.20 1.14e+06 ------1 E Calibration Curve : y = (-64389.797447) + (25516095.618234)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999195 Hept. epoxide : Single Peak Component Component Type : 15.485 min Retention Time Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area 0.0050 110981.00 31402.90 ------1 А 0.0100 219920.00 62911.46 ------ --в 1 С 0.0500 1239897.57 357220.78 ------ -----1 0.1000 2455174.68 704054.05 ------D 1 0.1500 3805678.03 1.09e+06 ----Е 1 Calibration Curve : y = (-32270.521957) + (25374615.531610)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999615 gamma chiordane Component Type : Single Peak Component Retention Time : 15.769 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates 0.0050 118257.40 32548.71 ---А _____ 1 В 228893.80 64946.68 ------ -----0.0100 1 0.0500 1291870.70 379501.28 ------ -----С 1 D 0.1000 2598359.70 762909.70 ------1 Е 0.1500 4085039.40 1.20e+06 ------ --1 Calibration Curve : y = (-49681.138416) + (27208973.625654)x + (0.000000)x^2 + (0.000000)x^3 : 0.999154 R-squared alpha chlordane Component Type : Single Peak Component Retention Time : 16.076 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
	A	0.0050	110705.40	31065.45		<u></u>	1
	C	0.0500	1199439.60	339743.57			1
	D	0.1000	2378312.50	673961.92			1
	E	0.1500	3733917.60	1.06e+06			1
	Calibration C R-squared	Curve : y : 0	e = (-37796.63).999224	86529) + (24	839303.5110	017)x + (0.00	0000)x^2 + (0.000000)x^3
4,	4'-DDE						
•	Component 1	Гуре	: Single Pe	eak Compon	ent		
	Retention Tir Search Wind	ne	: 16.262 m	າມກ 00_%			
	Reference C	omponer	nt :	.00 /0			
	Find peak clo	osest to e	expected RT i	n window			
	Calibrating A Curve will iar	rea versu	us Amount us prigin	ing a 1st Or	der Hit		
	Amounts will	not be s	caled prior to	the regressi	on		
	Weighting fac	ctor for th	ne regression:	: 1	0000/		
	Component s	standard	punty percen	tage : 100.0	000%		
	User Values						
	Label :	0 500000	h				
	Value 1 :	5.000000)				
	Value 3 :	0.000000)				
	Value 4 :	0.000000)				
	value 5.	0.000000	,				
	Calibration L	evel					
	Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
	A	0.0050	102179.96	30524.15			1
	B	0.0100	206062.76	62302.07			1
	C	0.0500	1250043.70	381709.07			1
	E	0.1500	3984153.60	1.22e+06			1
	Calibration C R-squared	urve:y	r = (-63853.19 .999243	9423) + (26	659285.1328	333)x + (0.00	0000)x^2 + (0.000000)x^3
-							
E	ndosulitan I Component 1	Evpe	: Single Pr	eak Compon	ent		
	Retention Tir	ne	: 16.385 m	nin			
	Search Wind	ow	: 5.00 s, 0.	.00 %			
	Find peak clo	omponer osest to e	expected RT i	n window			
	Calibrating A	rea versu	us Amount us	ing a 1st Or	der Fit		
	Curve will igr	nore the o	origin colod prior to	the regressi	~n		
	Weighting fai	ctor for th	e regression:	1	011		
	Component s	standard	purity percen	tage : 100.0	000%		
	l Iser Values						
	Label :						
	Value 1 : I	0.500000)				
	Value 2 :	0.000000	,)				
	Value 4 :	0.000000)				
	Value 5 :	0.000000)				
	Calibration Le	evel					
	Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
	Δ	0.0050	110157 04	30328 56			1
	B	0.0100	214396.94	59520.86			1
	ç	0.0500	1190087.30	333650.11			1
	D F	0.1000	2353209.20	062249.66 1 02e+06			1
	-	0.1000	0000021110	1.020100			•
	Calibration C	urve : y	= (-27711.38	9975) + (24	306143.2703	308)x + (0.00	0000)x^2 + (0.000000)x^3

: 0.999584 R-squared

Dieldrin Component Type : Single Peak Component Retention Time : 16.928 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area 0.0050 116591.70 31793.23 --1 A 0.0100 232249.40 64174.33 ----В 1 С 0.0500 1346525.80 378292.88 -----____ ____ 1 D 0.1000 2713923.80 755883.79 -----1 Е 0.1500 4246287.60 1.19e+06 ------1 Calibration Curve : y = (-55255.253370) + (28355092.828504)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999357 Endrin Component Type : Single Peak Component Retention Time : 17.447 min Search Window 5.00 s, 0.00 % • Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Height ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area 0.0050 120152.89 28808.87 ----А 1 в 0.0100 226480.90 58027.36 -----1 С 0.0500 1202358.40 329586.98 -----1 D 0.1000 2424613.64 659129.28 ------1 ε 0.1500 3748343.11 1.02e+06 -----1 Calibration Curve : y = (-27367.629171) + (24948529.543863)x + (0.000000)x² + (0.000000)x³ : 0.999600 R-squared 4.4'-DDD Component Type : Single Peak Component : 17.591 min Retention Time Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level ISTD Amt. ISTD Resp. # Replicates Level Name Amount Height Area 0.0050 90822.51 26377.21 ----1 А 183903.50 53096.79 ----- -----1 В 0.0100 С 0.0500 1078355.60 318520.93 ------ -----1 0.1000 2168968.26 640616.24 ------D 1 0.1500 3370849.49 981231.21 ------1 E Calibration Curve : y = (-41734.186130) + (22544666.744032)x + (0.000000)x^2 + (0.000000)x^3 : 0.999574 R-squared Endosulfan II Component Type : Single Peak Component : 17.943 min Retention Time Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates А 0.0050 109885.30 29117.37 -----1 0.0100 211709.40 56332.09 ------В 1 0.0500 1165848.30 315792.62 ------1 С D 0.1000 2268419.90 614873.72 ------1 Е 0.1500 3540898.00 957223.35 -----1 Calibration Curve : y = (-21806.222775) + (23510450.837696)x + (0.000000)x^2 + (0.000000)x^3 : 0.999366 R-squared 4,4'-DDT Component Type : Single Peak Component Retention Time : 18.203 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A	0.0050	90105.60	25382.30			1		
	B	0.0100	183517.00	51993.20			1		
	D	0.0500	2272435.60	662127.87			1		
	E	0.1500	3529670.20	1.02e+06			1		
	Calibration C R-squared	urve:y :0	= (-50266.63 .999607	1160) + (23	662487.7246	567)x + (0.00	0000)x^2 + (0.000000)x^	3	
E	ndrin aldehyd Component Retention Tir Search Wind Reference C Find peak dd Calibrating A Curve will ig Amounts will Weighting fa Component s User Values Label : Value 1 : Value 2 : Value 3 :	le Type me low omponen osest to e note the o not be se ctor for th standard 0.5000000 5.000000	: Single Pe : 18.868 m : 5.00 s, 0. t : expected RT i us Amount us prigin caled prior to ne regression: purity percen	eak Compon in .00 % ing a 1st Or the regressi 1 tage : 100.0	der Fit on 000%				
	Value 3 : Value 4 : Value 5 : Calibration L	0.000000 0.000000 0.000000)						
	Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A	0.0050	85758.80	22464.39			1		
	B	0.0100	162309.10	41967.97			1		
	D	0.0500	1691286.03	437137.87			1		
	E	0.1500	2640904.00	686949.94			1		
	Calibration C R-squared	Curve:y :0	= (-7460.562 .999198	:661) + (174	85838.8573	52)x + (0.000	000)x^2 + (0.000000)x^3	\$	
M	Methoxychlor Component Type : Single Peak Component Retention Time : 19.243 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1								
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000))))						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A	0.0050	56192.70	16317.10			1		
	ь С	0.0100	600219.60	171251.03			1		
	Ď	0.1000	1161118.40	332185.62			1		
	E	0.1500	1756625.80	505104.97			1		
	Calibration Curve : y = (-260.641034) + (11703679.698953)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999832								

Endo. Sulfate Component Type : Single Peak Component : 19.821 min Retention Time : 5.00 s, 0.00 % Search Window Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area Height 102886.60 26058.22 ----А 0.0050 1 197109.20 49781.59 ------ ----в 0.0100 1 С 0.0500 1073572.60 276673.38 ------1 0.1000 2072923.80 539455.73 ------1 D ---Е 0.1500 3246466.30 838321.01 ------1 Calibration Curve : y = (-16939.448789) + (21516367.441099)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999227 Endrin ketone Component Type : Single Peak Component Retention Time : 20.434 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates А 0.0050 112146.20 27459.34 ------ -----1 0.0100 219519.60 53898.74 ------ -----В 1 0.0500 1237282.00 307130.12 ------ -----С 1 D 0.1000 2425006.60 602555.78 ------1 Е 0.1500 3780653.20 941798.54 ------1 Calibration Curve : y = (-30049.583181) + (25158271.479058)x + (0.000000)x² + (0.000000)x³ R-squared : 0.999440

Software Version	: 6.2.1.0.104:0104	Date	: 07/24/2009 07:35:12
Reprocess Number	: buf1938: 102855		
Operator	: tchrom	Sample Name	: 9030585
Sample Number	:	Study	: ICAL 0.15
AutoSampler	: BUILT-IN	Rack/Vial	: 1/7
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 16:02:48	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42007.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42007.rst Inst Method : h:\turbo6\6890-06\6890-06\6890-66\62-process.mth from H:\TURBO6\6890-06\6-seq42\6A42007.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42007.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42007.rst Report Format File: h:\turbo6\6890-06\6890-06\65amp.rpt Sequence File : H:\TURBO6\6890-06\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.		Height [µV]
1	9.01	13667		в	0.01367	4247.48
2	11.22	4707018	alpha-BHC	В	0.15000	1.46e+06
3	12.08	4290829	gamma-BHC	В	0.15000	1.31e+06
4	12.32	1709671	beta-BHC	В	0.15000	513910.88
5	12.77	4419651	delta-BHC	В	0.15000	1.33e+06
6	12.94	16863		V	0.01686	3929.52
7	13.29	3948258	Heptachlor	В	0.15000	1.18e+06
8	14.02	3811422	Aldrin	В	0.15000	1.14e+06
9	14.74	3952		В	0.00395	1280.19
10	14.83	31646		V	0.03165	8791.96
11	15.27	15214		В	0.01521	4256.32
12	15.48	3805678	Hept, epoxide	V	0.15000	1.09e+06
13	15.76	4085039	gamma chlordane	В	0.15000	1.20e+06
14	16.07	3733918	alpha chlordane	В	0.15000	1.06e+06
15	16.25	3984154	4,4'-DDE	В	0.15000	1.22e+06
16	16.38	3650028	Endosulfan I	V	0.15000	1.02e+06
17	16.92	4246288	Dieldrin	В	0.15000	1.19e+06
18	17.20	51260		В	0.05126	13921.02
19	17.44	3748343	Endrin	8	0.15000	1.02e+06
20	17.58	3370849	4,4'-DDD	V	0.15000	981231.21
21	17.93	3540898	Endosulfan II	В	Ó.15000	957223.35
22	18.19	3529670	4,4'-DDT	в	0.15000	1.02e+06

07/24/2009 07:35:12 Result: H:\TURBO6\6890-06\6-seq42\6A42007.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
23	18.60	10279		в	0.01028	2877.74
24	18.72	2922		В	0.00292	1085.88
25	18.86	2640904	Endrin aldehyde	V	0.15000	686949.94
26	19.24	1756626	Methoxychlor	В	0.15000	505104.97
27	19.41	10387	•	В	0.01039	3101.42
28	19.81	3246466	Endo. Sulfate	В	0.15000	838321.01
29	20.10	36843		В	0.03684	9122.69
30	20.43	3780653	Endrin ketone	В	0.15000	941798.54
		72199396			3.19303	2.07e+07

Software Version	: 6.2.1.0.104:0104	Date	: 07/24/2009 07:35:15
Reprocess Number	: buf1938: 102856		
Operator	: tchrom	Sample Name	: 9030584
Sample Number	:	Study	: ICAL 0.10
AutoSampler	: BUILT-IN	Rack/Vial	: 1/8
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	; 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 16:38:40	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42008.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42008.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6A42008.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42008.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42008.rst Report Format File: h:\turbo6\6890-06\6a-007-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42008.rst

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.		Height [µV]
1	8.19	3535		в	0.00353	1317.31
2	9.01	10034		В	0.01003	3188.51
3	11.22	2951163	alpha-BHC	В	0.10000	927945.19
4	12.07	2698992	gamma-BHC	В	0.10000	827260.29
5	12.32	1086216	beta-BHC	В	0.10000	325971.18
6	12.76	2754822	delta-BHC	В	0.10000	838079.11
7	12.94	11142		Е	0.01114	2585.85
8	13.29	2539545	Heptachlor	В	0.10000	770765.95
9	14.02	2421194	Aldrin	В	0.10000	727111.12
10	14.82	17748		В	0.01775	5386.38
11	15.27	9591		в	0.00959	2729.44
12	15.48	2455175	Hept. epoxide	V	0.10000	704054.05
13	15.76	2598360	gamma chlordane	в	0.10000	762909.70
14	16.07	2378312	alpha chlordane	в	0.10000	673961.92
15	16.25	2535969	4,4'-DDE	в	0.10000	773533.60
16	16.38	2353209	Endosulfan I	V	0.10000	662249.66
17	16.92	2713924	Dieldrin	в	0.10000	755883.79
18	17.20	32935		В	0.03294	9057.44
19	17.44	24 2 4614	Endrin	в	0.10000	659129.28
20	17.58	2168968	4,4'-DDD	V	0.10000	640616.24
21	17.93	2268420	Endosulfan II	в	0.10000	614873.72
22	18.19	2272436	4,4'-DDT	в	0.10000	662127.87

07/24/2009 07:35:15 Result: H:\TURBO6\6890-06\6-seq42\6A42008.rst

Peak	Time	Area	Component	ΒĿ	NG	Height
#	[min]	[uV-sec]	Name		CONCENTRATION	[µV]
				·		
23	18.60	7003		в	0.00700	1976.80
24	18.72	3033		В	0.00303	1048.52
25	18.86	1691286	Endrin aldehyde	V	0.10000	437137.87
26	19.24	1161118	Methoxychlor	В	0.10000	332185.62
27	19.41	7640	·	в	0.00764	2283.13
28	19.81	2072924	Endo. Sulfate	В	0.10000	539455.73
29	20.10	22506		в	0.02251	5652.49
30	20.43	2425007	Endrin ketone	В	0.10000	602555.78
		<u>_</u>			<u> </u>	
		46096820			2.12517	1.33e+07

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102857	Date	: 07/24/2009 07:35:17
Operator	: tchrom	Sample Name	: 9071072
Sample Number	:	Study	: ICAL 0.05
AutoSampler	: BUILT-IN	Rack/Vial	: 1/9
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 17:14:41	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42009.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42009.rst Inst Method : h:\turbo6\6890-06\6890-06\6890-06\6890-06\6-seq42\6A42009.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42009.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42009.rst Report Format File: h:\turbo6\6890-06\6a-07-23-09.mth

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	9.02	5802		в	0.00580	1884.70
2	11.23	1422693	alpha-BHC	в	0.05000	454715.04
3	12.09	1313186	gamma-BHC	В	0.05000	403167.56
4	12.33	551639	beta-BHC	В	0.05000	164775.79
5	12.78	1331458	delta-BHC	В	0.05000	404736.00
6	13.30	1267229	Heptachlor	В	0.05000	386320.86
7	13.60	9672	•	В	0.00967	2706.14
8	14.03	1194589	Aldrin	В	0.05000	363472.70
9	14.83	8582		В	0.00858	2707.90
10	15.28	4387		В	0.00439	1348.12
11	15.48	1239898	Hept. epoxide	V	0.05000	357220.78
12	15.77	1291871	gamma chlordane	В	0.05000	379501.28
13	16.08	1199440	alpha chlordane	В	0.05000	339743.57
14	16.26	1250044	4,4'-DDE	В	0.05000	381709.07
15	16.39	1190087	Endosulfan I	V	0.05000	333650.11
16	16.93	1346526	Dieldrin	В	0.05000	378292.88
17	17.21	17384		В	0.01738	4766.48
18	17.45	1202358	Endrin	в	0.05000	329586.98
19	17.59	1078356	4,4'-DDD	V	0.05000	318520.93
20	17.94	1165848	Endosulfan II	в	0.05000	315792.62
21	18.20	1126622	4,4'-DDT	в	0.05000	326602.33
22	18.61	3953		В	0.00395	1171.02

07/24/2009 07:35:17 Result: H:\TURBO6\6890-06\6-seq42\6A42009.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
23	18.87	890478	Endrin aldehyde	B	0.05000	228942.24
24	19.24	600220	Methoxychlor	В	0.05000	171251.03
25	19.42	3562	·	в	0.00356	1153.85
26	19.82	1073573	Endo. Sulfate	В	0.05000	276673.38
27	20.11	11960		В	0.01196	3057.01
28	20.43	1237282	Endrin ketone	В	0.05000	307130.12
29	24.59	134110		В	0.13411	18670.84
		23172807			1,19941	6.66e+06

•

Software Version	: 6.2.1.0.104:0104	Date : 07/2-	4/2009 07:35:19
Reprocess Number	: buf1938: 102858		
Operator	: tchrom	Sample Name : 9071	083
Sample Number	:	Study : ICAL	. 0.01
AutoSampler	: BUILT-IN	Rack/Vial : 1/10	
Instrument Name	: HP6890-06	Channel : A	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.4	5 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject : 1500	0.00000
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	: 07/23/2009 17:50:35	Cycle : 4	
Raw Data File : H:\TU Result File : H:\TURB(Inst Method : h:\turbof	RBO6\6890-06\6-seq42\6A42010.ra D6\6890-06\6-seq42\6A42010.rst \\6890-06\6890-6 12-5-08-ins from	w <modified> </modified>	w

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42010.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42010.rst Report Format File: h:\turbo6\6890-06\6a-mp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	11.22	217038	alpha-BHC	в	0.01000	67803.66
2	12.08	208947	gamma-BHC	в	0.01000	62929.01
3	12.32	103633	beta-BHC	в	0.01000	30429.29
4	12.77	205191	delta-BHC	в	0.01000	61484.26
5	13.29	215244	Heptachlor	В	0.01000	64186.05
6	14.02	193159	Aldrin	В	0.01000	57786.96
7	15.48	219920	Hept. epoxide	В	0.01000	62911.46
8	15.76	228894	gamma chlordane	В	0.01000	64946.68
9	16.07	213023	alpha chlordane	В	0.01000	59073.96
10	16.26	206063	4,4'-DDE	В	0.01000	62302.07
11	16.38	214397	Endosulfan I	V	0.01000	59520.86
12	16.92	232249	Dieldrin	в	0.01000	64174.33
13	17.44	226481	Endrin	в	0.01000	58027.36
14	17.59	183903	4,4'-DDD	V	0.01000	53096.79
15	17.94	211709	Endosulfan II	в	0.01000	56332.09
16	18.20	183517	4,4'-DDT	в	0.01000	51993.20
17	18.86	162309	Endrin aldehyde	в	0.01000	41967.97
18	19.24	111199	Methoxychlor	в	0.01000	31680.86
19	19.81	197109	Endo. Sulfate	в	0.01000	49781.59
20	20.43	219520	Endrin ketone	В	0.01000	53898.74
					`	

^{0.20000 1.11}e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102859	Date	: 07/24/2009 07:35:21
Operator	: tchrom	Sample Name	: 9071084
Sample Number	:	Study	: ICAL 0.005
AutoSampler	: BUILT-IN	Rack/Vial	: 1/11
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 18:26:42	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42011.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42011.rst Inst Method : h:\turbo6\6890-06\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42011.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42011.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42011.rst Report Format File: h:\turbo6\6890-06\6a-07-23-09.mth from H:\TURBO6\6890-06\6-seq42\6A42011.rst

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	11.22	104357	alpha-BHC	в	0.00500	32529.05
2	12.07	102120	gamma-BHC	в	0.00500	30487.42
3	12.32	53416	beta-BHC	В	0.00500	15497.55
4	12.77	99325	delta-BHC	в	0.00500	29585.29
5	13.29	107993	Heptachlor	в	0.00500	31928.31
6	14.03	95257	Aldrin	в	0.00500	28243.18
7	15.48	110981	Hept. epoxide	в	0.00500	31402.90
8	15.76	118257	gamma chlordane	в	0.00500	32548.71
9	16.07	110705	alpha chlordane	в	0.00500	31065.45
10	16.26	102180	4,4'-DDE	В	0.00500	30524.15
11	16.38	110157	Endosulfan I	V	0.00500	30328.56
12	16.92	116592	Dieldrin	В	0.00500	31793.23
13	17.44	120153	Endrin	в	0.00500	28808.87
14	17.58	90823	4,4'-DDD	V	0.00500	26377.21
15	17.94	109885	Endosulfan II	В	0.00500	29117.37
16	18.20	90106	4.4'-DDT	В	0.00500	25382.30
17	18.86	85759	Endrin aldehyde	В	0.00500	22464.39
18	19.24	56193	Methoxychlor	В	0.00500	16317.10
19	19.82	102887	Endo. Sulfate	В	0.00500	26058.22
20	20.43	112146	Endrin ketone	В	0.00500	27459.34

0.10000 557918.60

Software Version	: 6.2.1.0.104:0104	Date	: 07/24/2009 07:40:06
Reprocess Number	: buf1938: 102860		
Operator	: tchrom	Sample Name	: 9030626
Sample Number	:	Study	: 2nd Source
AutoSampler	: BUILT-IN	Rack/Vial	: 1/12
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 19:02:28	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42012.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42012.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6A42012.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42012.rst Calib Method : h:\turbo6\6890-06\6a-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42012.rst Report Format File: h:\turbo6\6890-06\6890-06\60-42.seq Sequence File : H:\TURBO6\6890-06\66-42.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
11.23	BB	1447108	alpha-BHC	0.04928	459319.55	-1.4	11.18 -	11.28
12.08	BB	1340225	gamma-BHC	0.04969	412265.89	-0.6	12.03 -	12.13
12.33	BB	565527	beta-BHC	0.05105	167923.42	2.1	12.28 -	12.38
12.77	BB	1350410	delta-BHC	0.04910	415800.48	-1.8	12.72 -	12.82
13.29	BB	1286922	Heptachlor	0.05055	390396.14	1.1	13.24 -	13.34
14.03	BB	1227221	Aldrin	0.05062	371119.33	1.2	13.98 -	14.08
15.48	ΒB	1281121	Hept. epoxide	0.05176	367401.66	3.5	15.43 -	15.53
15.77	BB	1293324	gamma chlordane	0.04936	375244.63	-1.3	15.72 -	15.82
16.07	BB	1208173	alpha chlordane	0.05016	340782.20	0.3	16.02 -	16. 12
16.26	ΒV	1269851	4,4'-DDE	0.05003	391226.63	0.1	16.21 -	16.31
16.38	VB	1205074	Endosulfan I	0.05072	338831.29	1.4	16.33 -	16.43
16.93	BB	1370108	Dieldrin	0.05027	382368.74	0.5	16.88 -	16.98
17.44	ΒV	1227318	Endrin	0.05029	333249.86	0.6	17.39 -	17.49
17.59	VB	1094070	4,4'-DDD	0.05038	323123.68	0.8	17.54 -	17.64
17.94	BB	1175525	Endosulfan II	0.05093	318471.96	1.9	17.89 -	17.99
18.20	BB	1139567	4,4'-DDT	0.05028	332121.47	0.6	18.15 -	18.25
18.87	VB	974552	Endrin aldehyde	0.05616	255770.14	12.3	18.82 -	18.92
19.24	BB	598792	Methoxychlor	0.05118	171399.61	2.4	19.19 -	19.29
19.82	BB	1061100	Endo. Sulfate	0.05010	274243.87	0.2	19.77 -	19.87
20.43	BB	1234846	Endrin ketone	0.05028	305845.83	0.6	20.38 -	20.48
		23350835		1.01219	6.73e+06			

 TotalChrom Method File H:\TURBO6\6890-06\6B-(07-23-09).mth

 Printed by
 : BescoD on: 07/24/2009 07:26:11

 Created by
 : BescoD on: 07/24/2009 06:59:42

 Edited by
 : BescoD on: 07/24/2009 07:26:04

 Number of Times Edited
 : 4

 Number of Times Calibrated
 : 2636

 Description:
 CURVE 07-23-09

Global Sample Information

Default Sample Volume	: 1.000 ul
Quantitation Units	: ng
Void Time	: 0.000 min
Correct amounts during calibration	: Yes
Convert unknowns to concentration units	: Yes
Reject outliers during calibration	: No

An External Standard calibration will be used Unknown peaks will be quantitated using a response factor of 1.000000e+06 First peak will be relative retention reference

Component Information

 alpha-BHC (2C)

 Component Type
 : Single Peak Component

 Retention Time
 : 12.820 min

 Search Window
 : 5.00 s, 0.00 %

 Reference Component :
 :

 Find peak closest to expected RT in window

 Calibrating Area versus Amount using a 1st Order Fit

 Curve will ignore the origin

 Amounts will not be scaled prior to the regression

 Weighting factor for the regression: 1

 Component standard purity percentage : 100.0000%

Area

User Values

Label : Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level Level Name Amount

ISTD Amt. ISTD Resp. # Replicates

11111

A	0.0050	55058.60	16237.12		
B	0.0100	111801.20	33154.31	***********	
С	0.0500	692901.30	209761.45		
D	0.1000	1427291.60	432250.99	***********	
E	0.1500	2286305.00	681689.95		

Height

Calibration Curve : $y = (-48435.795610) + (15287417.467814)x + (0.000000)x^2 + (0.000000)x^3$ R-squared : 0.998362

gamma-BHC (2C)

Component Type: Single Peak ComponentRetention Time: 13.784 minSearch Window: 5.00 s, 0.00 %Reference Component :Find peak closest to expected RT in windowCalibrating Area versus Amount using a 1st Order FitCurve will ignore the originAmounts will not be scaled prior to the regressionWeighting factor for the regression: 1Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.020000

Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Processed by: <u>DYB</u> <u>1</u> 124,09	
Reviewed by: <u>NUT 8 24,09</u>	

Calibration Lev	/el .mount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A (B (C (D (E (0.0050 0.0100 0.0500 0.1000 0.1500	54029.20 106466.20 628264.80 1271798.20 2017688.80	15296.92 30354.75 181864.74 369343.87 586050.08			1 1 1 1 1	
Calibration Cur R-squared	rve:y :0.	= (-32259.70 998827	9534) + (13	3458874.918	281)x + (0.00	0000)x^2 + (().000000)x^3
beta-BHC (2C) Component Ty Retention Time Search Window Reference Cor Find peak clos Calibrating Are Curve will igno Amounts will n Weighting facts Component sta	rpe w nponen est to e va versu re the o ot be so or for th andard p	: Single Pe : 13.995 m : 5.00 s, 0 t : xpected RT i s Amount us rigin aled prior to e regression: purity percen	eak Compor in .00 % n window ing a 1st Or the regress : 1 tage : 100.0	nent rder Fit ion 0000%			
User Values Label : Value 1 : 0. Value 2 : 0. Value 3 : 0. Value 4 : 0. Value 5 : 0.	020000 000000 000000 000000 000000						
Calibration Lev Level Name A	/el mount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A (0) B (0) C (0) D (0) E (0)	0.0050 0.0100 0.0500 0.1000 0.1500	28927.60 55319.20 274557.50 525505.30 814939.40	8211.24 15428.84 76912.28 149557.24 232077.88			1 1 1 1	
Calibration Cur R-squared	rve:y :0.	= (1082.725 999429	536) + (537	7254.960082	2)x + (0.00000	00)x^2 + (0.00	00000)x^3
delta-BHC (2C) Component Type : Single Peak Component Retention Time : 14.752 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%							
User Values Label : Value 1 : 0. Value 2 : 0. Value 3 : 0. Value 4 : 0. Value 5 : 0.	020000 000000 000000 000000 000000						
Calibration Lev	vel Imount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A () B () C () D () E ()	0.0050 0.0100 0.0500 0.1000 0.1500	51371.00 101551.30 609220.40 1238174.80 1984350.40	14458.46 28653.74 174459.87 353178.73 567090.83			1 1 1 1 1	

Calibration Curve : y = (-36618.075345) + (13230978.468500)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.998354

Heptachlor (2C) Component Type : Single Peak Component : 14.944 min Retention Time Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area 0.0050 54132.60 15262.75 --1 А В 0.0100 103561.20 29129.76 ---1 С 569915.00 162908.51 ----0.0500 1 D 0.1000 1137804.80 327536.36 ------1 1 Ε Calibration Curve : $y = (-16834.226440) + (11839374.608405)x + (0.000000)x^{2} + (0.000000)x^{3}$ R-squared : 0.999319 Aldrin (2C) Component Type : Single Peak Component Retention Time 15.754 min Search Window 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Height ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area А 0.0050 46390.70 12929.16 ------1 В 0.0100 89690.80 24771.26 -----1 С 0.0500 509856.60 145746.10 ------ -----1 D 0.1000 1032204.80 293209.74 -----1 0.1500 1629985.70 464830.40 -----1 F Calibration Curve ; y = (-22501.175339) + (10859156.977685)x + (0.000000)x^2 + (0.000000)x^3 : 0.998963 R-squared Hept. epoxide (2C) Component Type : Single Peak Component : 17.110 min Retention Time Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000					
Calibration L Level Name	.evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	48056.60 91165.20 475920.60 941743.40 1464410.20	12821.79 24192.77 128536.19 257707.61 397231.88			1 1 1 1 1
Calibration C R-squared	Curve:y :0.	= (-7773.377 999469	291) + (971	4802.814130	6)x + (0.0000	00)x^2 + (0.000000)x^3
gamma chlordane (2C) Component Type : Single Peak Component Retention Time : 17.549 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%						
User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000					
Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	48722.60 90705.90 468633.80 930134.40 1458010.90	12867.95 23970.13 128488.28 258057.35 405024.76			1 1 1 1
Calibration C R-squared	Curve : y : 0.	= (-9056.291 .999178	950) + (965	5520.82460	7)x + (0.0000	00)x^2 + (0.000000)x^3
alpha chlordane (2C) Component Type : Single Peak Component Retention Time : 17.884 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.040000 Value 2 : 0.000000						
Value 3 : Value 4 : Value 5 :	0.000000 0.000000 0.000000					

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07/24/2009 07:26:11 Method: H:\TURBO6\6890-06\6B-(07-23-09).mth

C L	alibration Lo evel Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A		0.0050 0.0100	52574.33 93111.52	12358.20 22817.63			1 1	
C		0.0500	450266.81	121564.77		*****	1	
E		0.1000	1375657.70	376024.55			1	
C F	Calibration C R-squared	urve:y :0.	= (-199.0246 999292	66) + (9064	310.847060)	x + (0.00000	0)x^2 + (0.000000)x^3	
	dosulfan I (2 Component 1 Retention Tir Search Wind Reference Ca Tind peak clo Calibrating A Curve will igr mounts will Veighting fac Component s	C) fype ne omponent osest to e rea versu ore the o not be so ctor for th standard p	: Single Pe : 18.027 m : 5.00 s, 0. t : xpected RT in s Amount us rigin caled prior to e regression: purity percent	eak Compon in 00 % n window ing a 1st Ord the regressi 1 age : 100.00	ent der Fit on 000%			
ι	Jser Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000 0.000000						
C L	Calibration Lo	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A		0.0050	43146.87	11301.33			1	
E		0.0100	81984.68	21270.77	***********	****	1	
)	0.0500	424131.39 831714.07	222807.61			1	
E		0.1500	1292420.90	346955.62			1	
C F	Calibration C C-squared	urve:y :0.	= (-4830.583 999480	741) + (856	3653.410311	l)x + (0.0000	00)x^2 + (0.000000)x^3	
4,4 C F S F F C C A V C	4,4'-DDE (2C) Component Type : Single Peak Component Retention Time : 18.219 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%							
L	Jser Values Label : Value 1 : 1 Value 2 : 4 Value 3 : 0 Value 4 : 0 Value 5 : 0	0.500000 5.000000 0.000000 0.000000 0.000000 0.000000						
C L	alibration Le	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A	L	0.0050	39113.90	11123.79			1	
B		0.0100	76012.00	21421.05			1	
D)	0.1000	861972.00	253252.13			1	
E		0.1500	1359380.10	396104.31			1	

Calibration Curve : $y = (-17494.611158) + (9051517.637435)x + (0.000000)x^2 + (0.000000)x^3$ R-squared : 0.999022

Dieldrin (2C) Component Type : Single Peak Component Retention Time : 18.647 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area Height 11106.05 ----- --0.0050 42811.80 A В 0.0100 82756.60 21290.14 ------1 ¢ 0.0500 451139.00 119291.97 ----1 D 0.1000 902887.60 241590.63 ------ --1 Е 0.1500 1418012.20 378842.47 ------ ---1 Calibration Curve : y = (-14360.400046) + (9426695.529817)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999171 Endrin (2C) : Single Peak Component Component Type Retention Time : 19.337 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates А 0.0050 34827.33 8825.28 -----1 0.0100 68368.71 17102.75 -----В 1 365582.43 92776.64 -----С 0.0500 1 D 0.1000 731969.60 188776.12 ------1 0.1500 1140837.75 294594.42 ------1 Е Calibration Curve : y = (-9706.312366) + (7587673.951510)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999420 4,4'-DDD (2C) Component Type : Single Peak Component : 19.500 min Retention Time : 5.00 s, 0.00 % Search Window Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level ISTD Amt. ISTD Resp. # Replicates Height Level Name Amount Area 0.0050 32013.67 8907.43 -----1 А 62663.09 17234.02 -----1 в 0.0100 С 0.0500 346075.17 96045.83 -----1 693429.20 194480.44 -----D 0.1000 1 0.1500 1091229.55 305094.21 ----1 Ε Calibration Curve ; y = (-12232.467395) + (7258961.693510)x + (0.000000)x^2 + (0.000000)x^3 : 0.999115 R-squared Endosulfan II (2C) Component Type : Single Peak Component Retention Time : 19.793 min : 5.00 s, 0.00 % Search Window Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates 39654.20 1 A 0.0050 9873.61 ----- ---74934.60 18512.62 -----В 0.0100 1 383984.30 95912.53 -----1 С 0.0500 D 0.1000 739003.60 187421.46 ------ ---1 Е 1 Calibration Curve : y = (-430.453599) + (7569200.850785)x + (0.000000)x^2 + (0.000000)x^3 : 0.999499 R-squared 4,4'-DDT (2C) Component Type : Single Peak Component Retention Time : 20.200 min : 5.00 s, 0.00 % Search Window Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration	Level e Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	26168.80	7224.77			1
B	0.0100	53036.50	14386.39		*******	1
	0.0500	636640.80	175952.00			1
Ē	0.1500	990675.50	274922.61			1
Calibration R-squared	Curve : y : 0	r = (-13491.0 .999536	74017) + (6	631094.3991	199)x + (0.00	0000)x^2 + (0.000000)x^3
Endrin aldeh Componen Retention T Search Wir Reference Find peak of Calibrating Curve will i Amounts w Weighting f	yde (2C) t Type Time dow Componer closest to e Area versu gnore the c ill not be s factor for th t standard	: Single F : 20.524 r : 5.00 s, (at : expected RT us Amount us origin caled prior to ne regression purity percer	Peak Compo nin).00 % in window sing a 1st C o the regres: 1: 1 ntage : 100.	nent Irder Fit sion 0000%		
User Value Label Value 1 Value 2 Value 3 Value 4 Value 5	s : 0.500000 : 5.000000 : 0.000000 : 0.000000 : 0.000000					
Calibration Level Name	Level e Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	30800.00	7607.28			1
В	0.0100	57635.00	14045.24			1
C D	0.0500	288067.20	132193 27			1
Ē	0.1500	833217.62	205904.05			1
Calibration R-squared	Curve : y : 0	= (3974.252 9.999214	2925) + (548	4029.68886	4)x + (0.0000	000)x^2 + (0.000000)x^3
Endo. Sulfate Componen Retention T Search Wir Reference Find peak o Calibrating Curve will i Amounts w Weighting t Componen	e (2C) t Type Time dow Componer closest to e gnore the c gnore the c factor for th t standard	: Single F : 21.138 r : 5.00 s, (at : expected RT us Amount us origin caled prior to ne regression purity percer	Peak Compo nin).00 % in window sing a 1st C o the regres: n: 1 ntage : 100.	nent Irder Fit sion 0000%		
User Value Label Value 1 Value 2 Value 3 Value 4 Value 5	s 0.500000 5.000000 0.000000 0.0000000 0.0000000)))				
Calibration	Level e Amount	Area	Height	ISTD Amt	ISTD Resp	, #Replicates
A	0.0050	38791.80	9242.55	<u></u>		• 1
B	0.0100	72802.50	17121.97 87432.23	/		- 1 - 1
D	0.1000	685633.00	166109.94	,		• 1
Е	0.1500	1059884.40	259213.42	2		- 1
Calibration R-squared	Curve : y : 0	= (4487.498 .999332	3069) + (698	4292.72905	8)x + (0.0000	000)x^2 + (0.000000)x^3

Methoxychlor (2C) Component Type : Single Peak Component Retention Time : 21.669 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area Height 0.0050 16136.00 4126.55 -----А 1 _____ в 0.0100 32421.00 8154.20 ------ -----1 С 0.0500 177565.70 43531.81 -----1 D 0.1000 337505.40 83300.11 ------ -----1 0.1500 504323.60 125958.44 ------ -----E 1 Calibration Curve : y = (1660.904941) + (3363959.286649)x + (0.000000)x^2 + (0.000000)x^3 : 0.999554 R-squared Endrin ketone (2C) Component Type : Single Peak Component **Retention Time** : 22.451 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Height ISTD Amt. ISTD Resp. # Replicates Level Name Amount Area 40685.20 0.0050 8285.08 -----1 А В 0.0100 78100.20 15715.54 ------1 0.0500 С 414075.10 83906.04 -----1 D 0.1000 792447.60 161627.00 ------1 E 0.1500 1223581.75 251736.55 -----1 Calibration Curve : y = (-1025.638038) + (8107993.777019)x + (0.000000)x^2 + (0.000000)x^3

R-squared : 0.999532

Software Version	: 6.2.1.0.104:0104	Date	: 07/24/2009 07:23:51
Operator	: tchrom	Sample Name	: 9030585
Sample Number	•	Study	: ICAL 0.15
AutoSampler	: BUILT-IN	Rack/Vial	: 1/7
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.00000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 16:02:48	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42007.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42007.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42007.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42007.rst Callb Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42007.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.74	2951		В	0.00295	1006.04
2	10.47	9197		В	0.00920	2584.56
3	12.81	2286305	alpha-BHC (2C)	В	0.15000	681689.95
4	13.77	2017689	gamma-BHC (2C)	В	0.15000	586050.08
5	13.99	814939	beta-BHC (2C)	В	0.15000	232077.88
7	14.74	1984350	delta-BHC (2C)	В	0.15000	567090.83
8	14.94	1779818	Heptachlor (2C)	В	0.15000	515138.02
9	15.17	3972		В	0.00397	1078.27
10	15.75	1629986	Aldrin (2C)	в	0.15000	464830.40
11	16.70	13243		в	0.01324	3679.85
12	17.10	1464410	Hept. epoxide (2C)	В	0.15000	397231.88
13	17.54	1458011	gamma chlordane (2C)	В	0.15000	405024.76
14	17.88	1375658	alpha chlordane (2C)	₿	0.15000	376024.55
15	18.02	1292421	Endosulfan I (2C)	V	0.15000	346955.62
16	18.21	1359380	4,4'-DDE (2C)	В	0.15000	396104.31
17	18.64	1418012	Dieldrin (2C)	₿	0.15000	378842.47
19	19.33	1140838	Endrin (2C)	В	0.15000	294594.42
20	19.49	1091230	4,4'-DDD (2C)	V	0.15000	305094.21
21	19.79	1144569	Endosulfan II (2C)	В	0.15000	291547.98
22	20.19	990676	4,4'-DDT (2C)	В	0.15000	274922.61
23	20.52	833218	Endrin aldehyde (2C)	В	0.15000	205904.05
24	20.77	13528		V	0.01353	2557.28

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07/24/2009 07:23:51 Result: H:\TURBO6\6890-06\6-seq42\6b42007.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
25	21.13	1059884	Endo. Sulfate (2C)	В	0.15000	259213.42
26	21.66	504324	Methoxychlor (2C)	В	0.15000	125958.44
27	21.95	2966		В	0.00297	755.36
28	22.20	9554		В	0.00955	2009.40
29	22.44	1223582	Endrin ketone (2C)	V	0.15000	251736.55
		26924710			3.05541	7.37e+06

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Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102848	Date	: 07/24/2009 07:23:56
Operator	: tchrom	Sample Name	: 9030584
Sample Number	:	Study	: ICAL 0.10
AutoSampler	: BUILT-IN	Rack/Vial	: 1/8
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	; 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 16:38:40	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42008.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42008.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42008.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42008.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42008.rst Report Format File: h:\turbo6\6890-06\6b-07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42008.rst Report Format File: h:\turbo6\6890-06\6b-07-23-09

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.74	3311		в	0.00331	1142.06
2	10.47	7196		В	0.00720	1995.32
4	12.81	1427292	alpha-BHC (2C)	В	0.10000	432250.99
5	13.77	1271798	gamma-BHC (2C)	В	0.10000	369343.87
6	13.98	525505	beta-BHC (2C)	В	0.10000	149557.24
7	14.74	1238175	delta-BHC (2C)	В	0.10000	353178.73
8	14.93	1137805	Heptachlor (2C)	В	0.10000	327536.36
9	15.17	2461	,	В	0.00246	691.04
10	15.75	1032205	Aldrin (2C)	В	0.10000	293209.74
11	16.70	8397		в	0.00840	2310.35
12	17.10	941743	Hept. epoxide (2C)	В	0.10000	257707.61
13	17.54	930134	gamma chlordane (2C)	В	0.10000	258057.35
14	17.88	882653	alpha chlordane (2C)	В	0.10000	240252.20
15	18.02	831714	Endosulfan I (2C)	V	0.10000	222807.61
16	18.21	861972	4,4'-DDE (2C)	В	0.10000	253252.13
17	18.64	902888	Dieldrin (2C)	В	0.10000	241590.63
19	19.33	731970	Endrin (2C)	В	0.10000	188776.12
20	19.49	693429	4,4'-DDD (2C)	V	0.10000	194480.44
21	19.78	739004	Endosulfan II (2C)	В	0.10000	187421.46
22	20.19	636641	4,4'-DDT (2C)	в	0.10000	175952.00
23	20.52	537621	Endrin aldehyde (2C)	в	0.10000	132193.27
24	20.77	7460	,	в	0.00746	1633.35

07/24/2009 07:23:56 Result: H:\TURBO6\6890-06\6-seq42\6b42008.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
25	21.13	685633	Endo. Sulfate (2C)	В	0.10000	166109.94
26	21.66	337505	Methoxychlor (2C)	В	0.10000	83300.11
27	21.95	2129	• • • •	В	0.00213	560.11
28	22.20	5786		В	0.00579	1249.94
29	22.44	792448	Endrin ketone (2C)	В	0.10000	161627.00
		17174873			2.03674	4.70e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102850	Date	: 07/24/2009 07:24:01
Operator	: tchrom	Sample Name	: 9071072
Sample Number	:	Study	: ICAL 0.05
AutoSampler	: BUILT-IN	Rack/Vial	: 1/9
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 17:14:41	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42009.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42009.rst Inst Method : h:\turbo6\6890-06\6890-612-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42009.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42009.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42009.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.74	2505		в	0.00251	868.50
2	10.48	4581		В	0.00458	1227.98
4	12.82	692901	alpha-BHC (2C)	В	0.05000	209761.45
5	13.78	628265	gamma-BHC (2C)	В	0.05000	181864.74
6	13.99	274557	beta-BHC (2C)	В	0.05000	76912.28
7	14.75	609220	delta-BHC (2C)	B	0.05000	174459.87
8	14.94	569915	Heptachlor (2C)	в	0.05000	162908.51
9	15.75	509857	Aldrin (2C)	в	0.05000	145746.10
10	16.71	4373		в	0.00437	1209.12
11	17.11	475921	Hept. epoxide (2C)	в	0.05000	128536.19
12	17.55	468634	gamma chlordane (2C)	в	0.05000	128488.28
13	17.88	450267	alpha chlordane (2C)	в	0.05000	121564.77
14	18.03	424131	Endosulfan I (2C)	V	0.05000	112848.18
15	18.22	427277	4,4'-DDE (2C)	в	0.05000	124388.59
16	18.65	451139	Dieldrin (2C)	в	0.05000	119291.97
17	19.34	365582	Endrin (2C)	в	0.05000	92776.64
18	19.50	346075	4,4'-DDD (2C)	V	0.05000	96045.83
19	19.79	383984	Endosulfan II (2C)	В	0.05000	95912.53
20	20.20	314818	4,4'-DDT (2C)	в	0.05000	86468.35
21	20.52	288067	Endrin aldehyde (2C)	в	0.05000	70354.96
22	20.78	4426		в	0.00443	998.14
23	21.14	365378	Endo. Sulfate (2C)	В	0.05000	87432.23
, I

07/24/2009 07:24:01 Result: H:\TURBO6\6890-06\6-seq42\6b42009.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
24	21.67	177566	Methoxychlor (2C)	В	0.05000	43531.81
25	22.21	2109		B	0.00211	525.81
26	22.45	414075	Endrin ketone (2C)	В	0.05000	83906.04
27	29.05	19765		В	0.01976	2091.23
		8675389			1.03776	2.35e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102852	Date	: 07/24/2009 07:24:05
Operator	: tchrom	Sample Name	: 9071083
Sample Number	;	Study	: ICAL 0.01
AutoSampler	: BUILT-IN	Rack/Vial	: 1/10
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 17:50:35	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42010.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42010.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42010.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42010.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42010.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	12.81	111801	alpha-BHC (2C)	B	0.01000	33154.31
2	13.77	106466	gamma-BHC (2C)	B	0.01000	30354.75
3	13.98	55319	beta-BHC (2C)	B	0.01000	15428.84
4	14.74	101551	delta-BHC (2C)	B	0.01000	28653.74
5	14.93	103561	Heptachlor (2C)	B	0.01000	29129.76
6	15.75	89691	Aldrin (2C)	B	0.01000	24771.26
7	17.10	91165	Hept. epoxide (2C)	B	0.01000	24192.77
8	17.54	90706	gamma chlordane (2C)	B	0.01000	23970.13
9	17.88	93112	alpha chlordane (2C)	B	0.01000	22817.63
10	18.02	81985	Endosulfan I (2C)	V	0.01000	21270.77
11	18.21	76012	4,4'-DDE (2C)	8	0.01000	21421.05
12	18.64	82757	Dieldrin (2C)	8	0.01000	21290.14
13	19.33	68369	Endrin (2C)	8	0.01000	17102.75
14	19.50	62663	4,4'-DDD (2C)	V	0.01000	17234.02
15	19.79	74935	Endosulfan II (2C)	B	0.01000	18512.62
16	20.20	53037	4,4'-DDT (2C)	B	0.01000	14386.39
17	20.52	57635	Endrin aldehyde (2C)	B	0.01000	14045.24
18	21.14	72802	Endo. Sulfate (2C)	B	0.01000	17121.97
19	21.67	32421	Methoxychlor (2C)	B	0.01000	8154.20
20	22.45	78100	Endrin ketone (2C)	B	0.01000	15715.54

0.20000 418727.86

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102854	Date	: 07/24/2009 07:24:10
Operator	: tchrom	Sample Name	: 9071084
Sample Number	;	Study	: ICAL 0.005
AutoSampler	: BUILT-IN	Rack/Vial	; 1/11
Instrument Name	: HP6890-06	Channel	; B
Instrument Serial #	: CN10520010	A/D mV Range	; 1000
Delay Time	: 0.00 min	End Time	; 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 18:26:42	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42011.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42011.rst

Inst Method : h:\turbo6\6890-06\6890-06\6890-66 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42011.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42011.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42011.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name			Height [µV]
1	12.81	55059	alpha-BHC (2C)	В	0.00500	16237.12
2	13.78	54029	gamma-BHC (2C)	в	0.00500	15296.92
3	13.99	28928	beta-BHC (2C)	В	0.00500	8211.24
4	14.75	51371	delta-BHC (2C)	в	0.00500	14458.46
5	14.94	54133	Heptachlor (2C)	В	0.00500	15262.75
6	15.75	46391	Aldrin (2C)	В	0.00500	12929.16
7	17.11	48057	Hept. epoxide (2C)	В	0.00500	12821.79
8	17.55	48723	gamma chlordane (2C)	В	0.00500	12867.95
9	17.88	52574	alpha chlordane (2C)	В	0.00500	12358.20
10	18.02	43147	Endosulfan I (2C)	V	0.00500	11301.33
11	18.21	39114	4,4'-DDE (2C)	В	0.00500	11123.79
12	18.64	42812	Dieldrin (2C)	В	0.00500	11106.05
13	19.33	34827	Endrin (2C)	В	0.00500	8825.28
14	19.50	32014	4,4'-DDD (2C)	V	0.00500	8907.43
15	19.79	39654	Endosulfan II (2C)	в	0.00500	9873.61
16	20.20	26169	4,4'-DDT (2C)	в	0.00500	7224.77
17	20.52	30800	Endrin aldehvde (2C)	B	0.00500	7607.28
18	21.14	38792	Endo. Sulfate (2C)	в	0.00500	9242.55
19	21.67	16136	Methoxychlor (2C)	в	0.00500	4126.55
20	22.45	40685	Endrin ketone (2C)	В	0.00500	8285.08

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102862	Date	: 07/24/2009 07:42:20
Operator	: tchrom	Sample Name	: 9030626
Sample Number	:	Study	: 2nd Source
AutoSampler	: BUILT-IN	Rack/Vial	: 1/12
Instrument Name	: HP6890-06	Channel	; В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 19:02:28	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42012.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42012.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42012.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42012.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42012.rst Report Format File: h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42012.rst Report Format File: h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42012.rst

Sequence File : H:\TURBO6\6890-06\6D-42.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.81	BB	691199	alpha-BHC (2C)	0.04838	208398.69	-3.2	12.76 -	12.86
13.78	BB	645712	gamma-BHC (2C)	0.05037	187254.17	0.7	13.73 -	13.83
13.99	BB	284178	beta-BHC (2C)	0.05265	79576.40	5.3	13.94 -	14.04
14.75	BB	635328	delta-BHC (2C)	0.05079	183049.18	1.6	14.70 -	14.80
14.94	BB	588361	Heptachlor (2C)	0.05112	168844.47	2.2	14.89 -	14.99
15.75	BB	534981	Aldrin (2C)	0.05134	152224.53	2.7	15.70 -	15.80
17.11	BB	505521	Hept. epoxide (2C)	0.05284	136687.97	5.7	17.06 -	17.16
17.55	BB	483562	gamma chlordane (2C)	0.05102	132895.35	2.0	17.50 -	17.60
17.88	ΒV	473550	alpha chlordane (2C)	0.05227	128218.93	4.5	17.83 -	17.93
18.03	VB	441874	Endosulfan I (2C)	0.05216	117272.17	4.3	17.98 -	18.08
18.22	BB	447197	4,4'-DDE (2C)	0.05134	128817.88	2.7	18.17 -	18.27
18.65	BB	471146	Dieldrin (2C)	0.05150	124511.32	3.0	18.60 -	18.70
19.34	ΒV	384887	Endrin (2C)	0.05200	97940.09	4.0	19.29 -	19.39
19.50	VB	368786	4,4'-DDD (2C)	0.05249	102503.43	5.0	19.45 -	19.55
19.79	BB	398657	Endosulfan II (2C)	0.05273	99505.61	5.5	19.74 -	19.84
20.20	BB	322620	4,4'-DDT (2C)	0.05069	88350.53	1.4	20.15 -	20.25
20.53	BB	329841	Endrin aldehyde (2C)	0.05942	80268.10	18.8	20.48 -	20.58
21.14	BB	372688	Endo. Sulfate (2C)	0.05272	89106.83	5.4	21.09 -	21.19
21.67	BB	180550	Methoxychlor (2C)	0.05318	44181.16	6.4	21.62 -	21.72
22.45	BB	423733	Endrin ketone (2C)	0.05239	85376.55	4.8	22.40 -	22.50
		8984373		1.04138	2.43e+06			

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102870	Date	: 07/24/2009 09:03:31
Operator	: tchrom	Sample Name	: 9030626
Sample Number	:	Study	: 2nd Source
AutoSampler	: BUILT-IN	Rack/Vial	: 1/13
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/24/2009 07:56:26	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42013.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42013.rst Inst Method : h:\turbo6\6890-06\6890-06\65-process.mth from H:\TURBO6\6890-06\6-seq42\6b42013.raw Proc Method : h:\turbo6\6890-06\65-process.mth from H:\TURBO6\6890-06\6-seq42\6b42013.rst Calib Method : h:\turbo6\6890-06\6b-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42013.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.80	BB	616821	alpha-BHC (2C)	0.04352	188497.81	-13.0	12.75 -	12.85
13.76	BB	559808	gamma-BHC (2C)	0.04399	166234.43	-12.0	13.71 -	13.81
13.97	BB	238119	beta-BHC (2C)	0.04408	68466.07	-11.8	13.92 -	14.02
14.73	BB	550916	delta-BHC (2C)	0.04441	160159.44	-1 1.2	14.68 -	14.78
14.92	BB	534073	Heptachlor (2C)	0.04653	155736.51	-6.9	14.87 -	14.97
15.73	BB	482121	Aldrin (2C)	0.04647	139114.58	-7.1	15.68 -	15.78
17.09	BB	461130	Hept. epoxide (2C)	0.04827	127154.95	-3.5	17.04 -	17.14
17.53	BB	435947	gamma chlordane (2C)	0.04609	120838.80	-7.8	17.48 -	17.58
17.87	ΒV	426961	alpha chlordane (2C)	0.04713	117332.66	-5.7	17.82 -	17.92
18.01	VB	399968	Endosulfan I (2C)	0.04727	108720.71	-5.5	17.96 -	18.06
18.20	BB	406326	4,4'-DDE (2C)	0.04682	119979.50	-6.4	18.15 -	18.25
18.63	BB	429353	Dieldrin (2C)	0.04707	115006.25	-5.9	18.58 -	18.68
19.32	ΒV	359862	Endrin (2C)	0.04871	92453.91	-2.6	19.27 -	19.37
19.49	VB	341596	4,4'-DDD (2C)	0.04874	96073.40	-2.5	19.44 -	19.54
19.78	BB	378134	Endosulfan II (2C)	0.05001	95357.97	0.0	19.73 -	19.83
20.19	BB	328795	4,4'-DDT (2C)	0.05162	92090.67	3.2	20.14 -	20.24
20.51	BB	300613	Endrin aldehyde (2C)	0.05409	76002.56	8.2	20.46 -	20.56
21.12	BB	350591	Endo. Sulfate (2C)	0.04955	85279.74	-0.9	21.07 -	21.17
21.65	BB	182993	Methoxychlor (2C)	0.05390	46113.92	7.8	21.60 -	21.70
22.43	BB	406365	Endrin ketone (2C)	0.05025	83806.06	0.5	22.38 -	22.48
		8190493		0.95852	2.25e+06			

 TotalChrom Method File H:\TURBO6\6890-06\6a-SURR-(07-23-09).mth

 Printed by
 : BescoD on: 07/24/2009 07:15:34

 Created by
 : BescoD on: 07/24/2009 07:04:10

 Edited by
 : BescoD on: 07/24/2009 07:15:29

 Number of Times Edited
 : 7

 Number of Times Calibrated
 : 2621

 Description:
 SURR CURVE 07-23-09

Global Sample Information

Default Sample Volume	:	1.000 ul
Quantitation Units	:	ng
Void Time	:	0.000 min
Correct amounts during calibration	:	Yes
Convert unknowns to concentration units	:	Yes
Reject outliers during calibration	:	No

An External Standard calibration will be used Unknown peaks will be quantitated using a response factor of 1.000000e+06

First peak will be relative retention reference

Component Information

 Tetrachloro-m-xylene

 Component Type
 : Single Peak Component

 Retention Time
 : 9.632 min

 Search Window
 : 7.00 s, 0.50 %

 Reference Component :
 :

 Find peak closest to expected RT in window

 Calibrating Area versus Amount using a 1st Order Fit

 Curve will ignore the origin

 Amounts will not be scaled prior to the regression

 Weighting factor for the regression: 1

 Component standard purity percentage : 100.0000%

Area

User Values

Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level Level Name Amount

ISTD Amt. ISTD Resp. # Replicates

A	0.0050	74829.60	22864.07	 	
В	0.0100	152166.80	46100.50	 	
С	0.0500	807581.00	252735.28	 	
D	0.0750	1208116.30	378938.71	 	
E	0.1000	1752098.20	548862.85	 	

Height

Calibration Curve : y = (-30831.080684) + (17287280.430906)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.996334

Decachlorobiphenyl

Component Type: Single Peak ComponentRetention Time: 22.684 minSearch Window: 7.00 s, 0.50 %Reference Component :Find peak closest to expected RT in windowCalibrating Area versus Amount using a 1st Order FitCurve will ignore the originAmounts will not be scaled prior to the regressionWeighting factor for the regression: 1Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000

Value 5 : 0.000000

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Processed by: <u>1943</u> 7 124109	
Reviewed by: AV & DA 107	

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07/24/2009 07:15:34 Method: H:\TURBO6\6890-06\6a-SURR-(07-23-09).mth

Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	118976.00	26164.22			1
В	0.0100	245628.40	51886.43			1
С	0.0500	1106049.50	239588.13			1
D	0.0750	1746547.80	376371.55		***	1
Е	0.1000	2227042.20	494712.11			1

Calibration Curve : $y = (13839.661664) + (22396023.298663)x + (0.000000)x^2 + (0.000000)x^3$ R-squared : 0.998705

Software Version	: 6.2.1.0.104:0104	Date : 07/24/2009 07:13:5	57
Reprocess Number	: buf1938: 102835		
Operator	: tchrom	Sample Name : 9030661	
Sample Number	:	Study : ICAL 0.10	
AutoSampler	: BUILT-IN	Rack/Vial : 1/2	
Instrument Name	: HP6890-06	Channel : A	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.45 min	
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject : 1500.000000	
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	: 07/23/2009 13:03:00	Cycle : 1	

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42002.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42002.rst Inst Method : h:\turbo6\6890-06\6890-66\2-5-08-ins from H:\TURBO6\6890-06\6-seq42\6A42002.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42002.rst Calib Method : h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42002.rst Report Format File: h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42002.rst Report Format File: h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42002.rst Sequence File : H:\TURBO6\6890-06\65amp.rpt



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	7.37	3136		в	0.00314	1094.15
2	9.64	1752098	Tetrachloro-m-xylene	B	0.10000	548862.85
4	10.75	18238		B	0.01824	5591.00
5	12.52	9614		В	0.00961	2923.58
6	13.58	38283		в	0.03828	10317.32
7	15.73	9538		В	0.00954	2623.47
8	18.19	11009		в	0.01101	402.10
9	18.72	6576		в	0.00658	1767.16
10	22.22	3932		в	0.00393	1044.76
11	22.69	2227042	Decachlorobiphenyl	в	0.10000	494712.11
		4079467			0.30033	1.07e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102837	Date	: 07/24/2009 07:14:02
Operator	: tchrom	Sample Name	: 9030660
Sample Number	•	Study	: ICAL 0.075
AutoSampler	: BUILT-IN	Rack/Vial	: 1/3
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 13:39:00	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42003.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42003.rst Inst Method : h:\turbo6\6890-06\6890-66\2-5-08-ins from H:\TURBO6\6890-06\6-seq42\6A42003.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42003.rst Calib Method : h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42003.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	9.63	1208116	Tetrachloro-m-xylene	в	0.07500	378938.71
3	10.74	13444		в	0.01344	4103.94
4	12.51	6580		в	0.00658	2062.19
5	15.73	15479		В	0.01548	4187.31
6	16.12	2841		В	0.00284	819.46
7	18.39	26732		В	0.02673	325.99
8	21.68	28403		В	0.02840	204.19
9	22.69	1746548	Decachlorobiphenyl	В	0.07500	376371.55
		3048142			0 24348	767013 34

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102839	Date	: 07/24/2009 07:14:07
Operator	: tchrom	Sample Name	: 9061646
Sample Number	:	Study	: ICAL 0.05
AutoSampler	: BUILT-IN	Rack/Vial	; 1/4
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.00000 uf	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 14:14:53	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42004.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42004.rst Inst Method : h:\turbo6\6890-06\6890-06\6890-06\6890-06\6-seq42\6A42004.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42004.rst Calib Method : h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42004.rst Report Format File: h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42004.rst Sequence File : H:\TURBO6\6890-06\66samp.rpt



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.18	3047		В	0.00305	1107.56
2	9.63	807581	Tetrachloro-m-xylene	в	0.05000	252735.28
3	10.39	1873	•	в	0.00187	471.89
4	10.74	9051		в	0.00905	2821.70
5	12.50	3893		в	0.00389	1328.68
6	15.73	8178		в	0.00818	2329.27
7	16.11	1854		в	0.00185	607.24
8	18.39	20521		В	0.02052	297.69
9	18.72	4096		в	0.00410	1294.73
10	22.68	1106049	Decachlorobiphenyl	В	0.05000	239588.13
		1000444			0.45054	
		1900144			0.15251	002082.17

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102841	Date	: 07/24/2009 07:14:11
Operator	: tchrom	Sample Name	: 9071081
Sample Number	•	Study	: ICAL 0.01
AutoSampler	: BUILT-IN	Rack/Vial	: 1/5
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 14:50:54	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-seq42\6A42005.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6A42005.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6A42005.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42005.rst Calib Method : h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42005.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	9.64	152167	Tetrachloro-m-xviene	в	0.01000	46100.50
2	15.73	7296	· · · · · · · · · · · · · · · · · · ·	в	0.00730	2034.73
3	17.40	15649		в	0.01565	5161.87
4	22.69	245628	Decachlorobiphenyl	в	0.01000	51886.43
		420741			0.04295	105183.53

Software Version	:	6.2.1.0.104:0104	Date	:	07/24/2009 07:14:15
Reprocess Number	:	buf1938: 102843			
Operator	:	tchrom	Sample Name	:	9071082
Sample Number	:		Study	:	ICAL 0.005
AutoSampler	:	BUILT-IN	Rack/Vial	:	1/6
Instrument Name	:	HP6890-06	Channel	:	A
Instrument Serial #	:	CN10520010	A/D mV Range	::	1000
Delay Time	:	0.00 min	End Time	:	29.46 min
Sampling Rate	:	5.0000 pts/s			
Sample Volume	:	1.000000 ul	Area Reject	:	1500.000000
Sample Amount	:	1.0000	Dilution Factor	:	1.00
Data Acquisition Time	:	07/23/2009 15:26:47	Cycle	:	5
Raw Data File : H:\TUI	RE	3O6\6890-06\6-seq42\6A42006.raw <modified></modified>			
Result File : H:\TURBO	26	\6890-06\6-seq42\6A42006.rst			
Inst Method : h:\turbo6	6/6	890-06\6890-6 12-5-08-ins from H:\TURBO6\689	0-06\6-seq42\6A4	120	006.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-seq42\6A42006.rst Calib Method : h:\turbo6\6890-06\6a-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6A42006.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	9.64	74830	Tetrachloro-m-xylene	в	0.00500	22864.07
2	15.73	5929		в	0.00593	1658.15
3	17.40	15641		в	0.01564	5109.06
4	19.44	3596		в	0.00360	868.36
5	22.69	118976	Decachlorobiphenyl	в	0.00500	26164.22
		218971			0.03517	56663.86

TotalChrom Method File H:\TU	RI	BO6\6890)-06	6B-SURR-(0	7-23-09).mth <modified></modified>	
Printed by	:	BescoD	on:	07/24/2009	07:16:20	
Created by	:	BescoD	on:	07/24/2009	07:05:59	
Edited by	:	BescoD	on:	07/24/2009	07:16:20	
Number of Times Edited	:	3				
Number of Times Calibrated	:	2651				
Description: SURR CURVE	07	-23-09				
•						

Global Sample Information

Default Sample Volume	:	1.000 ul
Quantitation Units	:	ng
Void Time	:	0.000 min
Correct amounts during calibration	:	Yes
Convert unknowns to concentration units	:	Yes
Reject outliers during calibration	:	No

An External Standard calibration will be used Unknown peaks will be quantitated using a response factor of 1.000000e+06 First peak will be relative retention reference

Component Information

 Tetrachloro-m-xylene (2C)

 Component Type
 : Single Peak Component

 Retention Time
 : 11.023 min

 Search Window
 : 7.00 s, 0.50 %

 Reference Component :
 :

 Find peak closest to expected RT in window

 Calibrating Area versus Amount using a 1st Order Fit

 Curve will ignore the origin

 Amounts will not be scaled prior to the regression

 Weighting factor for the regression: 1

 Component standard purity percentage : 100.0000%

Area

User Values

Label	:	
Value 1	:	0.500000
Value 2	:	5.000000
Value 3	:	0.000000
Value 4	:	0.000000
Value 5	:	0.000000

Calibration Level Level Name Amount

ISTD Amt. ISTD Resp. # Replicates

11111

А	0.0050	47433.20	13846.07	 	
B	0.0100	95838.10	27711.09	 	
С	0.0500	464074.20	137169.79	 	
D	0.0750	692870.40	207011.55	 	
E	0.1000	1012138.70	302459.69	 	

Height

Calibration Curve : $y = (-12432.280238) + (9893816.671620)x + (0.000000)x^{2} + (0.000000)x^{3}$ R-squared : 0.995296

Decachlorobiphenyl (2C) Component Type : Single Peak Component Retention Time : 26.109 min Search Window : 7.00 s, 0.50 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1

Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000

Value 5 : 0.000000

1	
Proce	essed by: 1945 1124 109
Dovid	moden NY 8,20, AS
Kevit	wed by: <u>1990</u> <u>0 107107</u>

07/24/2009 07:16:20 Method: H:\TURBO6\6890-06\6B-SURR-(07-23-09).mth

Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	44969.20	6972.26			1
В	0.0100	86907.60	13227.46			1
С	0.0500	369957.00	56001.63			1
D	0.0750	554229.60	84198.18			1
E	0.1000	719192.50	110780.41			1

Calibration Curve : $y = (13659.497623) + (7112326.716196)x + (0.000000)x^{2} + (0.000000)x^{3}$ R-squared : 0.999687

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102836	Date	: 07/24/2009 07:14:00
Operator	: tchrom	Sample Name	: 9030661
Sample Number	:	Study	: ICAL 0.10
AutoSampler	: BUILT-IN	Rack/Vial	: 1/2
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 13:03:00	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42002.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42002.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42002.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42002.rst Calib Method : h:\turbo6\6890-06\6b-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42002.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.43	3108		В	0.00311	699.05
2	8.73	2740		В	0.00274	983.38
3	11.03	1012139	Tetrachloro-m-xylene (В	0.10000	302459.69
4	12.12	10146		В	0.01015	2946.19
5	14.28	5097		В	0.00510	1403.47
6	15.84	2422		В	0.00242	694.12
7	17.54	9666		В	0.00967	1533.48
8	19.82	3201		В	0.00320	645.64
9	20.41	2478		В	0.00248	732.12
10	26.11	719193	Decachlorobiphenyl (2C	в	0.10000	110780.41
		1770189			0.23886	422877.54

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102838	Date	: 07/24/2009 07:14:05
Operator	: tchrom	Sample Name	: 9030660
Sample Number	:	Study	: ICAL 0.075
AutoSampler	: BUILT-IN	Rack/Vial	: 1/3
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.00000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 13:39:00	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42003.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42003.rst Inst Method : h:\turbo6\6890-06\6890-66\69-process.mth from H:\TURBO6\6890-06\6-seq42\6b42003.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42003.rst Calib Method : h:\turbo6\6890-06\6b-seurr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42003.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

.

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
2	8.73	2676		в	0.00268	965.72
3	11.02	692870	Tetrachloro-m-xylene (в	0.07500	207011.55
4	12.12	7461		в	0.00746	2141.62
5	14.28	3496		в	0.00350	1017.25
6	15.84	2523		в	0.00252	771.06
8	17.56	4010		в	0.00401	1068.61
9	19.22	3514		В	0.00351	163.79
10	24.19	2145		в	0.00214	154.71
11	26.11	554230	Decachlorobiphenyl (2C	в	0.07500	84198.18
		1272925			0.17583	297492.51

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102840	Date	: 07/24/2009 07:14:09
Operator	: tchrom	Sample Name	: 9061646
Sample Number	:	Study	: ICAL 0.05
AutoSampler	: BUILT-IN	Rack/Vial	: 1/4
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 14:14:53	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42004.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42004.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42004.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42004.rst Calib Method : h:\turbo6\6890-06\6b-seur-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42004.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.73	2932		в	0.00293	1017.71
3	11.02	464074	Tetrachloro-m-xylene (В	0.05000	137169.79
4	12.12	5112	•	В	0.00511	1468.32
5	14.28	2183		В	0.00218	666.50
6	15.84	1804		В	0.00180	550.83
7	17.09	2037		В	0.00204	110.97
8	17.56	1585		В	0.00158	483.31
10	26.11	369957	Decachlorobiphenyl (2C	В	0.05000	56001.63
		849684			0.11565	197469.07

Software Version	: 6.2.1.0.104:0104	Date	: 07/24/2009 07:14:13
Reprocess Number	: buf1938: 102842		
Operator	: tchrom	Sample Name	: 9071081
Sample Number	:	Study	: ICAL 0.01
AutoSampler	: BUILT-IN	Rack/Vial	: 1/5
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 14:50:54	Cycle	: 4
		•	

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42005.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42005.rst Inst Method : h:\turbo6\6890-06\6890-66\690-66\690-06\6-seq42\6b42005.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42005.rst Calib Method : h:\turbo6\6890-06\6b-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42005.rst Report Format File: h:\turbo6\6890-06\6b-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42005.rst Sequence File : H:\TURBO6\6890-06\6b-q2_seq

Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	11.03	95838	Tetrachloro-m-xylene (в	0.01000	27711.09
2	17.56	2427	• · ·	в	0.00243	662.29
3	17.82	5206		В	0.00521	1726.79
4	26.12	86908	Decachlorobiphenyl (2C	В	0.01000	13227.46
		190378			0.02763	43327.63

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf1938: 102844	Date	: 07/24/2009 07:14:16
Operator	: tchrom	Sample Name	: 9071082
Sample Number		Study	: ICAL 0.005
AutoSampler	: BUILT-IN	Rack/Vial	: 1/6
Instrument Name	: HP6890-06	Channel	:В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 07/23/2009 15:26:47	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-seq42\6b42006.raw <Modified> Result File : H:\TURBO6\6890-06\6-seq42\6b42006.rst Inst Method : h:\turbo6\6890-06\6890-66\690-612-5-08-ins from H:\TURBO6\6890-06\6-seq42\6b42006.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-seq42\6b42006.rst Calib Method : h:\turbo6\6890-06\6b-surr-(07-23-09).mth from H:\TURBO6\6890-06\6-seq42\6b42006.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6D-42.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	11.03	47433	Tetrachloro-m-xylene (в	0.00500	13846.07
2	17.56	1899	•	в	0.00190	531.79
3	17.82	5468		в	0.00547	1748.99
4	26.13	44969	Decachlorobiphenyl (2C	в	0.00500	6972.26
						<u> </u>
		99769			0.01737	23099.12

TotalChrom Method File H:\TURBO6\6890-()6\6a-(09-28-09).mth	
Printed by : NearvM o	n: 09/28/2009 14:38:20	
Created by NearyM o	n [•] 09/28/2009 14·26·29	
Edited by : NearyM o	$n = \frac{100}{200} \frac{1000}{1400} \frac{112000}{1400}$	- 100 mark 200 mil 100 mil 200
Number of Times Edited : 2		
Number of Times Calibrated 2000	a survey of the first and the	
Number of Times Calibrated : 2000		6.0
Description: CURVE 07-23-09	(1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1, 1	8 101
	MAN 1	NUMBER OF THE OFFICE OF
Global Sample Information	Processed of the OEP	2 9 2009
Default Sample Volume	: 1.000 ul (D) VEL (
Quantitation Units	: ng	and the second s
Void Time	: 0.000 min Reviewed of	יין איז
Correct amounts during calibration	: Yes	
Convert unknowns to concentration units	: Yes	
Reject outliers during calibration	: No	
An External Standard calibration will be us Unknown peaks will be quantitated using a First peak will be relative retention referen	ed a response factor of 1.000000e+06 ce	
O manage Male ymatic y		
(alpha-BHC)		
Determinent Type : Single Peak Col	nponent	
Retention Time : 10.743 min		
Search Window : 5.00 S, 0.00 %		
Reference Component :		
Find peak closest to expected RT in windo		
Calibrating Area versus Amount using a 1	st Order Fit	
Curve will ignore the origin		
Amounts will not be scaled prior to the reg	ression	
weighting factor for the regression: 1	00.00000/	
Component standard purity percentage : 1	00.0000%	
User values		
Label :		
Value 1 : 0.020000		
Value 2 : 0.000000		
Value 3 : 0.000000		
Value 4 : 0.000000		
Value 5 : 0.000000		
Calibration Level	ht ICTD And ICTD Deen # Device tes	
Level Name Amount Area Heig	nt ISTD Amt. ISTD Resp. # Replicates	
A 0.0050 151027.00 4017		
B 0.0100 31/227.00 4917	5.18 1	
C 0.0500 4033570 80 63040		
D 0.0000 1955070.00 05910		
E 0.1000 5797 158.90 1.896	+00 1	
Calibration Curve : $y = (41182, 426075)$.	(30405020 784135)	
R-squared 0 000108	(39403929.104133)X + (0.000000)X 2 + (0.000000)X 3	
11-3quared . 0.353150		
gamma-BHC		
	nponent	
Retention Time : 11 590 min	··· - · · · · ·	
Search Window 5.00 s. 0.00 %		
Reference Component		
Find peak closest to expected RT in wind		
Calibrating Area vareue Amount using a 1	at Order Fitz	
Curve will ignore the origin		
Amounts will not be scaled prior to the rea	rection	
Meighting factor for the regression: 4	10001011	
Component standard purity percentage 1		
	00.0000 /0	

User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.020000 0.000000 0.000000 0.000000 0.000000								
Calibration L	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates			
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	145445.10 296382.80 1755205.80 3616315.24 5333639.79	45737.04 93617.40 559673.93 1.15e+06 1.68e+06			1 1 1 1			
Calibration Curve : y = (-41853.472536) + (36051605.394861)x + (0.000000)x^2 + (0.000000)x^3 R-squared									
Deta-BHC Component Type : Single Peak Component Retention Time : 11.839 min Search Window : 5.00 s, 0.00 % Reference Component :									
User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.020000 0.000000 0.000000 0.000000 0.000000								
Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	•		
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	73550.60 141567.40 726066.60 1444746.50 2116347.50	22619.96 43178.96 224381.24 449159.83 666554.68			1 1 1 1			
Calibration (R-squared	Curve : y	= (8354.355) 999717	233) + (1416	0338.62256	7)x + (0.0000	000)x^2 + (0.0	00000)x^3		
delta-BHC Component Retention Ti Search Wind Reference C Find peak cl Calibrating A Curve will ig Amounts wil Weighting fa Component	Type me dow Component osest to e: Area versu nore the o I not be sc octor for th standard p	: Single Pe : 12.280 m : 3.00 s, 0. t : xpected RT i s Amount us rigin aled prior to e regression: purity percen	eak Compon in .00 % ing a 1st Or the regressi : 1 tage : 100.0	der Fib on 000%					

User Values Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 **Calibration Level** Level Name Amount ISTD Amt. ISTD Resp. # Replicates Area Height A 0.0050 137282.00 42738.23 -В 0.0100 285536.20 89319.43 --1 С 0.0500 1761227.40 554518.52 -1 D 0.1000 3654456.60 1.14e+06 -------1 Ε 0.1500 5494271.60 1.73e+06 ------ -----1 Calibration Curve : y = (-72133.889898) + (37122041.535057)x + (0.000000)x² + (0.000000)x³ : 0.999930 R-squared_ Heptachlor Component Type : Single Peak Component Retention Time 12.799 min Search Window 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label Value 1 : 0.020000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area 0.0050 151737.70 47866.25 -----A B C 1 0.0100 302559.80 93908.45 -----1 0.0500 1690219.90 536977.86 ------1 D 0.1000 3381776.20 1.06e+06 ------1 1.55e+06 -----Е 0.1500 4979244.00 1 Calibration Curve : $y = (-7869.57)(182) + (33475825.799284)x + (0.000000)x^2 + (0.000000)x^3$ R-squared : 0.999756 Aldrin Component Type : Single Peak Component Retention Time : 13.529 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit. Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

	User Values Label : Value 1 : (Value 2 : (Value 3 : (Value 4 : (Value 5 : (0.020000 0.000000 0.000000 0.000000 0.000000						
	Calibration Lo Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 (0.1500 4	140038.40 281975.90 1624794.40 3278401.50 4898481.40	43087.90 87658.92 512449.76 1.02e+06 1.51e+06			1 1 1 1 1	
	Calibration C R-squared	urve : y = : 0.9	= (-30301.27 999954	8076) + (32	937136.2008	832)x + (0.00	0000)x^2 + (0	0.000000)x^3
	Retention Tin Search Wind Reference Co Find peak clo Calibrating A Curve will ign Amounts will Weighting fac Component s User Values Label : Value 1 : 0 Value 2 : 1 Value 3 : 0	ne ow pomponent isest to ex- rea versus iore the or not be sca tor for the tandard p 0.500000 5.000000 0.000000	: 14.977 m : 5.00 s, 0. : pected RT ir s Amount usi igin aled prior to e regression: urity percent	in 00 % Window ing a 1st Ord the regressi 1 age : 100.0	der Fit on 000%			
: : :	Value 4 : (Value 5 : ().000000						
	Calibration Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	147629.60 289518.20 1569598.86 3113537.02 1602260.41	43518.43 85644.75 470002.47 919427.90 1.36e+06			1 1 1 1 1	
- 	Calibration C B-squared		= (1380.3636 999822	65)) + (3084	3308.79459	5)x + (0.0000)00)x^2 + (0.0)00000)x^3
Ø	amma chłorda Component T Retention Tin Search Wind Reference Co Find peak clo Calibrating A Curve will ign Amounts will Weighting fac Component s) ane by omponent sest to ex rea versus ore the or not be sca tor for the tandard p	: Single Pe : 15.261 m : 5.00 s, 0.1 : pected RT ir s Amount usi igin aled prior to to regression: urity percent	eak Compon in 00 % ng a 1st Or the regressi 1 age : 100.0	ent der Fit on 000%			

	User Values Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000) } }								
	Calibration Level Level Name Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates				
	Å 0.0050 B 0.0100 C 0.0500 D 0.1000 E 0.1500	145980.00 288979.20 1615121.00 3283497.60 4909801.30	43693.14 87530.95 495210.78 1.00e+06 1.49e+06			1 1 1 1 1				
	Calibration Curve : $y = (-28769.027062) + (32975315.032723)x + (0.000000)x^2 + (0.000000)x^3 R-squared (0.999967)$									
a	alpha chlordane : Single Peak Component Component Type : 15.566 min Search Window : 3.00 s, 0.00 % Reference Component : : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%									
	User Values Label : Value 1 : 0.040000 Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000)))								
	Calibration Level Level Name Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates				
	A 0.0050 B 0.0100 C 0.0500 D 0.1000 E 0.1500	142514.60 279403.60 1523815.00 3076903.20 4558923.90	42214.35 84764.31 465248.08 943079.17 1.39e+06			1 1 1 1 1				
	Calibration Curve : y R-squared : 0	= (-11570.27 .999899)219) + (30	601305.7620	659)x + (0.00	0000)x^2 + (0	.000000)x^3			
4	,4'-DDE Component Type Retention Time Search Window Reference Componer Find peak closest to e Calibrating Area versu Curve will ignore the o Amounts will not be so Weighting factor for th Component standard	: Single Pe : 15.768 m : 3.00 s, 0.1 : sxpected RT ir as Amount usion origin caled prior to the regression: purity percent	ak Compon in 00 % Ng a 1st Ord the regressi 1 age : 100.00	ent der Fit on 000%						

User Value Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	s 0.500000 5.000000 0.000000 0.000000 0.000000						
Calibration	Level Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	128243.14 260026.67 1525925.73 3122164.58 4658355.67	39966.76 82307.46 484682.70 991819.37 1.49e+06			1 1 1 1	
Calibration R-squared	Curve : 0.	= (-39044.4 9 999936	2463) + (31	396628.433	003)x + (0.00	0000)x^2 + (0	1.000000)x^3
Endosulfan I Component Retention T Search Win Reference Find peak o Calibrating Curve will ig Amounts w Weighting f	Type ime dow Component losest to e: Area versu nore the o Il not be sc actor for the standard p	: Single Pe : 15.868 m : 5.00 s, 0. t : s Amount us rigin aled prior to e regression: purity percent	eak Compon in .00 % ing a 1st Or the regressi : 1 tage : 100.0	der Fit on 000%			
User Value: Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	3 0.500000 5.000000 0.000000 0.000000 0.000000 0.000000						
Calibration	Level Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	132452.16 260436.63 1409819.87 2817111.12 4162708.73	38168.51 75402.71 415621.47 825242.18 1.21e+06			1 1 1 1	
Calibration R-squared	Curve : y : 0.	= (-2281.733 999846	833) + (279	17260.87328	80)x + (0.000	000)x^2 + (0.0	00000)x^3
Dieldrin Component Retention T Search Win Reference (Find peak o Calibrating Curve will ig Amounts wi Weighting f Component	Type ime dow Component losest to ex Area versu nore the o l not be sc actor for the standard p	: Single Pe : 16.409 m : 5.00 s, 0. : : s Amountus rigin aled prior to e regression: ourity percent	eak Compon in 00 % n window ing a 1st Or the regressi 1 tage : 100.0	ent der Fit on 000%			

09/28/2009 14:38:20 Method: H:\TURBO6\6890-06\6a-(09-28-09).mth

	User Values Label : Value 1 : 0 Value 2 : 0 Value 3 : 0 Value 4 : 0 Value 5 : 0	.040000 .000000 .000000 .000000 .000000							
	Calibration Le Level Name A A B C D E	vel mount 0.0050 0.0100 0.0500 0.1000 0.1500	Area 135109.20 271044.70 1555819.00 3163929.70 4708399.20	Height 39092.64 78524.15 459600.91 925794.81 1.37e+06	ISTD Amt.	ISTD Resp.	# Replicates 1 1 1 1 1 1		
	Calibration Cu R-squared	Irve : 7	≤ (-2 9797.58 999925	6218) + (31	692982.117(009)x + (0.00	0000)x^2 + (0	.000000)x^3	
E	Endrin Component Type : Single Peak Component Retention Time : 16.919 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%								
	User Values Label : Value 1 : 0 Value 2 : 0 Value 3 : 0 Value 4 : 0 Value 5 : 0	.040000 .000000 .000000 .000000 .000000							
	Calibration Le	vel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	112550.40 229729.30 1333091.00 2722950.40 4093839.80	31875.08 65488.63 385212.56 779503.57 1.18e+06			1 1 1 1 1		
	Calibration Cu R-squared	inte: y	= (-36869)81 999974	0148) + (27	544467.0926	646)x + (0.00	0000)x^2 + (0	0.000000)x^3	
R-squared : 0.999974 4,4'-DDD Component Type : Single Peak Component Retention Time : 17.085 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%									

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	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.040000 0.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	97094.00 198635.40 1158307.70 2367060.60 3507582.00	29402.80 61402.53 360039.17 729653.08 1.08e+06			1 1 1 1	
_	Calibration C B-equared	Curve y	= (-24839.73 .999868	6843) + (23	659930.515	250)x + (0.00	0000)x^2 + (0	.000000)x^3
E	ndosulfan II Component Retention Tir Search Wind Reference C Find peak cle Calibrating A Curve will igr Amounts will Weighting fa Component	Type me low omponen osest to e vrea versu nore the o not be so ctor for th standard j	: Single Pe : 17.412 m : 5.00 s, 0. t : xpected RT i is Amount us origin caled prior to e regression purity percen	eak Compor nin .00 % int a 1st Or the regress : 1 tage : 100.0	der Fit fon 000%			
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000	· · ·					
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	117407.40 234206.60 1246739.00 2510042.75 3732096.64	32009.06 64281.98 349709.30 706638.17 1.05e+06			1 1 1 1 1	
	Calibration C R-squared	prve : y : 0	= (-7453.197 .999948	637) + (250	08756.7502	27)x + (0.000	000)x^2 + (0.0	00000)x^3
4	,4'-DDT Component Retention Til Search Wind Reference C Find peak cle Calibrating A Curve will ign Amounts will Weighting fa Component) Type me low omponen osest to e vrea versu nore the c not be so ctor for th standard	: Single Pe : 17.692 m : 5.00 s, 0 t : xpected RT i s Amount us origin caled prior to e regression purity percen	eak Compor in .00 % ing a 1st Or the re gress : 1 tage : 100.0	der Fit			

User Values Label : Value 1 : 0.04000 Value 2 : 0.00000 Value 3 : 0.00000 Value 4 : 0.00000 Value 5 : 0.00000	0 0 0 0									
Calibration Level Level Name Amoun	t Area	Height	ISTD Amt.	ISTD Resp.	# Replicates					
A 0.0050 B 0.0100 C 0.0500 D 0.1000 E 0.1500	88502.30 185451.40 1135856.60 2369050.84 3661241.05	26623.16 56361.65 347787.30 724755.23 1.13e+06			1 1 1 1					
Calibration Curve R-squared	Calibration Curve (-64115.952628) + (24637085.178599)x + (0.000000)x^2 + (0.000000)x^3 R-squared 0.999600									
Perinn aldehyde Component Type: Single Peak Component extension TimeRetention Time: 18.332 min 5.00 s, 0.00 %Search Window: 5.00 s, 0.00 %Reference Component :Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%										
User Values Label : Value 1 : 0.50000 Value 2 : 5.00000 Value 3 : 0.00000 Value 4 : 0.00000 Value 5 : 0.00000	0 0 0 0 0									
Calibration Level	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates					
A 0.0050 B 0.0100 C 0.0500 D 0.1000 E 0.1500	78763.20 152870.40 828797.60 1669486.33 2464869.20	21484.91 41810.98 222024.79 448834.89 666706.82			1 1 1 1					
Calibration Curve : R-squared :	y = (-3222.005 0.999850	(14) + (165	42529.3941	70)x + (0.000	000)x^2 + (0.0)00000)x^3				
Methoxychlor Component Type Retention Time Search Window Reference Compone Find peak closest to Calibrating Area vers Curve will ignore the Amounts will not be a Weighting factor for the Component standard	: Single Pe : 18.739 m : 5.00 s, 0. nt : expected RT / us Amount us origin scaled prior to he regression: I purity percen	eak Compon in .00 % ing a 1st Or the regressi : 1 tage : 100.0	ent der Fit on 000%							

	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	49951.60 103943.20 558910.20 1132935.61 1729716.14	15130.56 31083.35 166648.56 339868.10 521778.92			1 1 1 1	
Calibration Curve y = (-13580, 78690) + (11566224.258905)x + (0.000000)x^2 + (0.000000)x R-squared : 0.999882								
¢	Endo. Sulfate Component Type : Single Peak Component Retention Time : 19.277 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%							
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	•
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	70186.00 151171.10 885395.10 1860879.00 2914743.80	18795.21 40461.99 240471.47 510023.61 791648.51			1 1 1 1 1	
	Calibration C R-squared	Curve : y	= (-55662.91 . <u>9991</u> 18	1976) + (19	557738.2853	340)x + (0.00	0000)x^2 + (0.9	000000)x^3
Endrin ketone Component Type : Single Peak Component Retention Time : 19.882 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%								

User Values Label : Value 1 : 0.500000

Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name	Amount	Area	Height	ISTD Amt.	ISID Resp.	# Replicates
				···-··	·	
A	0.0050	104000.60	26733.54			1
В	0.0100	211115.40	54440.55			1
С	0.0500	1170892.50	306246.38			1
D	0.1000	2395548.10	633840.29		***********	1
Ε	0.1500	3751262.64	990184.29			1

Calibration Curve : $y = (-50713.618188) + (25036150.260749)x + (0.000000)x^2 + (0.000000)x^3$ R-squared (0.0999231)

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98758	Date	: 09/28/2009 14:35:18
Operator	: tchrom	Sample Name	: 9090862-CCV1
Sample Number	: 0.15	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/15
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	:5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 11:30:02	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46215.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46215.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46215.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46215.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46215.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.50	2240		В	0.00224	900.98
2	8.58	21316		В	0.02132	7385.51
3	9.37	4677		В	0.00468	1548.77
4	10.77	5797159	alpha-BHC	В	0.15000	1.89e+06
5	11.50	3193		В	0.00319	1246.07
6	11.62	5333640	gamma-BHC	V	0.15000	1.68e+06
7	11.87	2116347	beta-BHC	В	0.15000	666554.68
8	12.31	5494272	delta-BHC	В	0.15000	1.73e+06
9	12.43	8455		В	0.00845	3347.34
10	12.83	4979244	Heptachlor	В	0.15000	1.55e+06
11	13.55	4898481	Aldrin	В	0.15000	1.51e+06
12	14.26	3751		В	0.00375	1386.81
13	14.35	42892		V	0.04289	12250.03
14	14.80	19872		В	0.01987	5746.46
15	15.00	4602260	Hept. epoxide	V	0.15000	1.36e+06
16	15.28	4909801	gamma chlordane	В	0.15000	1.49e+06

09/28/2009 14:35:18 Result: H:\TURBO6\6890-06\6-SEQ46\6A46215.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.	NG CONCENTRATION	Height [µV]
17	15.59	4558924	alpha chlordane	В	0.15000	1.39e+06
18	15.79	4658356	4,4'-DDE	В	0.15000	1.49e+06
19	15.89	4162709	Endosulfan I	V	0.15000	1.21e+06
20	16.43	4708399	Dieldrin	В	0.15000	1.37e+06
21	16.73	51413		В	0.05141	15277.50
22	16.94	4093840	Endrin	В	0.15000	1.18e+06
23	17.11	3507582	4,4'-DDD	В	0.15000	1.08e+06
24	17.27	10043		В	0.01004	2087.51
25	17.44	3732097	Endosulfan II	V	0.15000	1.05e+06
26	17.60	6087		В	0.00609	2151.58
27	17.71	3661241	4,4'-DDT	V	0.15000	1.13e+06
28	18.09	52368		В	0.05237	14367.81
29	18.35	2464869	Endrin aldehyde	В	0.15000	666706.82
30	18.61	4109	-	В	0.00411	1139.79
31	18.76	1729716	Methoxychlor	V	0.15000	521778.92
32	18.94	9333	•	V	0.00933	3042.05
33	19.30	2914744	Endo. Sulfate	В	0.15000	791648.51
34	19.58	21374		В	0.02137	5258.99
35	19.75	2992		В	0.00299	1049.33
36	19.90	3751263	Endrin ketone	V	0.15000	990184.29
37	20.62	7110		В	0.00711	777.98
38	22.09	19970		В	0.01997	2625.60
		00000400				0.4007

82366139

3.29120 2.48e+07

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98760	Date	: 09/28/2009 14:35:25
Operator	: tchrom	Sample Name	: 9090863-CCV1
Sample Number	: 0.10	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/16
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 12:05:55	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46216.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46216.rst

Inst Method : h:\turbo6\6890-06\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46216.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46216.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46216.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.48	3425		В	0.00343	1296.50
2	7.75	3061		В	0.00306	1167.65
3	8.56	29912		В	0.02991	10053.00
4	9.35	8855		В	0.00886	2905.32
5	10.75	4010065	alpha-BHC	В	0.10000	1.31e+06
6	11.47	3217		В	0.00322	1165.22
7	11.59	3616315	gamma-BHC	V	0.10000	1.15e+06
8	11.84	1444746	beta-BHC	В	0.10000	449159.83
9	12.28	3654457	delta-BHC	В	0.10000	1.14e+06
10	12.41	35999		E	0.03600	6909.31
11	12.80	3381776	Heptachlor	В	0.10000	1.06e+06
12	13.53	3278402	Aldrin	В	0.10000	1.02e+06
13	14.24	2870		В	0.00287	1029.27
14	14.32	30806		V	0.03081	8718.71
15	14.78	14547		В	0.01455	4120.80
16	14.98	3113537	Hept. epoxide	V	0.10000	919427.90

09/28/2009 14:35:25 Result: H:\TURBO6\6890-06\6-SEQ46\6A46216.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	15.26	3283498	gamma chlordane	В	0.10000	1.00e+06
18	15.57	3076903	alpha chlordane	В	0.10000	943079.17
19	15.77	3122165	4,4'-DDE	В	0.10000	991819.37
20	15.87	2817111	Endosulfan I	V	0.10000	825242.18
21	16.41	3163930	Dieldrin	В	0.10000	925794.81
22	16.71	33632		В	0.03363	10086.43
23	16.92	2722950	Endrin	В	0.10000	779503.57
.24	17.09	2367061	4,4'-DDD	В	0.10000	729653.08
25	17.25	6785		В	0.00679	1525.98
26	17.41	2510043	Endosulfan II	V	0.10000	706638.17
27	17.58	5783		В	0.00578	1981.55
28	17.69	2369051	4,4'-DDT	V	0.10000	724755.23
29	18.07	116767		В	0.11677	29361.23
30	18.34	1669486	Endrin aldehyde	V	0.10000	448834.89
31	18.60	3945		В	0.00394	1069.78
32	18.74	1132936	Methoxychlor	V	0.10000	339868.10
33	18.92	6679		В	0.00668	2112.49
34	19.28	1860879	Endo. Sulfate	В	0.10000	510023.61
35	19.56	15892		В	0.01589	3925.21
36	19.89	2395548	Endrin ketone	В	0.10000	633840.29
37	21.85	71684		В	0.07168	247.76
38	22.08	8341		В	0.00834	799.18

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2.40220 1.67e+07

Software Version	: 6.2.1.0.104:0104	Date	: 09/28/2009 14:35:32
Reprocess Number	: buf2048: 98762		
Operator	: tchrom	Sample Name	: 9091289-CCV1
Sample Number	: 0.05	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/17
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reiect	; 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 12:41:53	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46217.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46217.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46217.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46217.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46217.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	2745		В	0.00274	1068.57
2	8.56	19527		В	0.01953	6574.25
3	9.35	5460		В	0.00546	1810.40
4	10.75	1933571	alpha-BHC	В	0.05000	639100.91
5	11.59	1755206	gamma-BHC	В	0.05000	559673.93
6	11.84	726067	beta-BHC	В	0.05000	224381.24
7	12.28	1761227	delta-BHC	В	0.05000	554518.52
8	12.41	20352		Е	0.02035	4123.81
9	12.80	1690220	Heptachlor	В	0.05000	536977.86
10	13.53	1624794	Aldrin	В	0.05000	512449.76
11	14.32	14996		В	0.01500	4391.65
12	14.78	7477		В	0.00748	2134.51
13	14.98	1569599	Hept. epoxide	V	0.05000	470002.47
14	15.26	1615121	gamma chlordane	В	0.05000	495210.78
15	15.57	1523815	alpha chlordane	В	0.05000	465248.08
16	15.77	1525926	4,4'-DDE	В	0.05000	484682.70

09/28/2009 14:35:32 Result: H:\TURBO6\6890-06\6-SEQ46\6A46217.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	15.87	1409820	Endosulfan I	v	0.05000	415621.47
18	16.41	1555819	Dieldrin	В	0.05000	459600.91
19	16.71	16847		В	0.01685	5057.11
20	16.92	1333091	Endrin	В	0.05000	385212.56
21	17.09	1158308	4,4'-DDD	В	0.05000	360039.17
22	17.41	1246739	Endosulfan II	В	0.05000	349709.30
23	17.70	1135857	4,4'-DDT	В	0.05000	347787.30
24	18.07	69032		В	0.06903	17923.35
25	18.34	828798	Endrin aldehyde	В	0.05000	222024.79
26	18.74	558910	Methoxychlor	В	0.05000	166648.56
27	19.28	885395	Endo. Sulfate	В	0.05000	240471.47
28	19.56	6365		В	0.00636	1763.97
29	19.89	1170893	Endrin ketone	В	0.05000	306246.38
30	22.08	5449		В	0.00545	369.28
		27177425			1.16825	8.24e+06
Software Version	: 6.2.1.0.104:0104 · buf2048· 98764	Date	: 09/28/2009 14:35:39			
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Operator	: tchrom	Sample Name	: 9091538-CCV1			
Sample Number	: 0.01	Study	: PEST ICAL			
AutoSampler	: BUILT-IN	Rack/Vial	: 1/18			
Instrument Name	: HP6890-06	Channel	: A			
Instrument Serial #	: CN10520010	A/D mV Range	: 1000			
Delay Time	: 0.00 min	End Time	: 29.47 min			
Sampling Rate	: 5.0000 pts/s					
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000			
Sample Amount	: 1.0000	Dilution Factor	: 1.00			
Data Acquisition Time	: 09/28/2009 13:17:47	Cycle	: 4			

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46218.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46218.rst Inst Method : h:\turbo6\6890-06\6890-612-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46218.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46218.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46218.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL.	NG CONCENTRATION	Height [µV]
1	8.56	6345		В	0.00635	2196.02
2	10.75	314234	alpha-BHC	В	0.01000	102805.18
3	11.60	296383	gamma-BHC	В	0.01000	93617.40
4	11.85	141567	beta-BHC	В	0.01000	43178.96
5	12.29	285536	delta-BHC	В	0.01000	89319.43
6	12.81	302560	Heptachlor	В	0.01000	93908.45
7	13.54	281976	Aldrin	В	0.01000	87658.92
8	14.98	289518	Hept. epoxide	В	0.01000	85644.75
9	15.27	288979	gamma chlordane	В	0.01000	87530.95
10	15.57	279404	alpha chlordane	В	0.01000	84764.31
11	15.78	260027	4,4'-DDE	В	0.01000	82307.46
12	15.88	260437	Endosulfan I	V	0.01000	75402.71
13	16.42	271045	Dieldrin	В	0.01000	78524.15
14	16.93	229729	Endrin	В	0.01000	65488.63
15	17.10	198635	4,4'-DDD	В	0.01000	61402.53
16	17.42	234207	Endosulfan II	В	0.01000	64281.98

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09/28/2009 14:35:39 Result: H:\TURBO6\6890-06\6-SEQ46\6A46218.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	17.70	185451	4,4'-DDT	В	0.01000	56361.65
18	18.08	17910		В	0.01791	4693.24
19	18.34	152870	Endrin aldehyde	В	0.01000	41810.98
20	18.75	103943	Methoxychlor	В	0.01000	31083.35
21	19.29	151171	Endo. Sulfate	В	0.01000	40461.99
22	19.89	211115	Endrin ketone	В	0.01000	54440.55
		4763043			0.22426	1.43e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98766	Date	: 09/28/2009 14:35:45
Operator	: tchrom	Sample Name	: 9091537-CCV1
Sample Number	: 0.005	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/19
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 13:53:55	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46219.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46219.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46219.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46219.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46219.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq

500 400 Response [mV] 300 200 SPL1 40.0 10.75 13,60 12.29 16.42 7.43 7.43 12.81 13.54 -19.89 18.895 23 (0) h 100 6 -8,56 ģ Q, ENDRIN ШŢ 1 Î GEMAR DELTA--HEPTA-ENDRIN-ALDRIN ENDO. ALPHA. internationality i 6 10 8 12 à 14 16 18 20 22 24 26 28 Time [min]

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.56	4005		В	0.00401	1417.43
2	10.75	151927	alpha-BHC	В	0.00500	49170.41
3	11.60	145445	gamma-BHC	В	0.00500	45737.04
4	11.85	73551	beta-BHC	В	0.00500	22619.96
5	12.29	137282	delta-BHC	В	0.00500	42738.23
6	12.81	151738	Heptachlor	В	0.00500	47866.25
7	13.54	140038	Aldrin	В	0.00500	43087.90
8	14.99	147630	Hept. epoxide	В	0.00500	43518.43
9	15.27	145980	gamma chlordane	В	0.00500	43693.14
10	15.58	142515	alpha chlordane	В	0.00500	42214.35
11	15.78	128243	4,4'-DDE	В	0.00500	39966.76
12	15.88	132452	Endosulfan I	V	0.00500	38168.51
13	16.42	135109	Dieldrin	В	0.00500	39092.64
14	16.93	112550	Endrin	В	0.00500	31875.08
15	17.10	97094	4,4'-DDD	В	0.00500	29402.80
16	17.43	117407	Endosulfan II	В	0.00500	32009.06

09/28/2009 14:35:45 Result: H:\TURBO6\6890-06\6-SEQ46\6A46219.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
17	17.70	88502	4,4'-DDT	В	0.00500	26623.16
18	18.09	11263		В	0.01126	2972.54
19	18.35	78763	Endrin aldehyde	В	0.00500	21484.91
20	18.75	49952	Methoxychlor	В	0.00500	15130.56
21	19.29	70186	Endo. Sulfate	В	0.00500	18795.21
22	19.89	104001	Endrin ketone	В	0.00500	26733.54
		2365633			0.11527	704317.92

Software Version	: 6.2.1.0.104:0104	Date	:	09/28/2009 14:59:54
Reprocess Number	: buf2048: 98768			
Operator	: tchrom	Sample Name	:	9091536
Sample Number	: 0.05	Study	:	ALT PEST
AutoSampler	: BUILT-IN	Rack/Vial	:	1/20
Instrument Name	: HP6890-06	Channel	:	Α
Instrument Serial #	: CN10520010	A/D mV Range	:	1000
Delay Time	: 0.00 min	End Time	:	29.48 min
Sampling Rate	: 5.0000 pts/s			
Sample Volume	: 1.000000 ul	Area Reject	:	3000.000000
Sample Amount	: 1.0000	Dilution Factor	:	1.00
Data Acquisition Time	: 09/28/2009 14:29:42	Cycle	:	1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46220.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46220.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46220.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46220.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46220.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL.	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
10.75	BB	2048447	alpha-BHC	0.05303	677304.60	6.1	10.70 -	10.80
11.60	BB	1864685	gamma-BHC	0.05288	594374.64	5.8	11.55 -	11.65
11.84	BB	770593	beta-BHC	0.05383	238784.39	7.7	11.79 -	11.89
12.28	ΒE	1877137	delta-BHC	0.05251	588469.71	5.0	12.23 -	12.33
12.80	BB	1801470	Heptachlor	0.05405	564866.62	8.1	12.75 -	12.85
13.53	BB	1770429	Aldrin	0.05467	552725.66	9.3	13.48 -	13.58
14.98	BB	1685104	Hept. epoxide	0.05459	504711.83	9.2	14.93 -	15.03
15.26	BΒ	1681298	gamma chlordane	0.05186	513401.10	3.7	15.21 -	15.31
15.57	BB	1606855	alpha chlordane	0.05289	498119.26	5.8	15.52 -	15.62
15.77	ΒV	1619278	4,4'-DDE	0.05282	515996.29	5.6	15.72 -	15.82
15.87	VB	1493267	Endosulfan I	0.05357	439323.82	7.1	15.82 -	15.92
16.41	BB	1646672	Dieldrin	0.05290	481838.13	5.8	16.36 -	16.46
16.92	BB	1413603	Endrin	0.05266	408600.87	5.3	16.87 -	16.97
17.09	BB	1242990	4,4'-DDD	0.05359	390903.37	7.2	17.04 -	17.14
17.42	BB	1341752	Endosulfan II	0.05395	372843.75	7.9	17.37 -	17.47
17.70	BB	1254290	4,4'-DDT	0.05351	382614.96	7.0	17.65 -	17.75

09/28/2009 14:59:54 Result: H:\TURBO6\6890-06\6-SEQ46\6A46220.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - Relative
18.34 18.75 19.28 19.89	VB BB BB BB	1065443 590147 931451 1237887	Endrin aldehyde Methoxychlor Endo. Sulfate Endrin ketone	0.06460 0.05220 0.05047 0.05147	290496.99 177700.00 252139.69 325558.62	29.2 4.4 0.9 2.9	18.29 - 18.39 18.70 - 18.80 19.23 - 19.33 19.84 - 19.94
28942799 1.07204 8.77e+06 Missing Component Report							ie INDB G-
All components were found							EAldelyde)

SEP 2 9 2009

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98782	Date	: 09/28/2009 15:39:15
Operator	: tchrom	Sample Name	: 9040971
Sample Number	:	Study	: INDBM MIX
AutoSampler	: BUILT-IN	Rack/Vial	: 1/21
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 15:09:19	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46221.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46221.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46221.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46221.rst Calib Method : h:\turbo6\6890-06\6a-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46221.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



$\begin{array}{cccccccccccccccccccccccccccccccccccc$	Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]		
$\begin{array}{cccccccccccccccccccccccccccccccccccc$									
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	1	7.77	3498		В	0.00350	1335.09		
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	2	8.57	8655		В	0.00866	3009.91		·
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	3	9.19	487306		В	0.48731	158598.23		
$\begin{array}{cccccccccccccccccccccccccccccccccccc$	4	11.86	274299	beta-BHC	В	0.01878	84101.52		
$ \begin{array}{cccccccccccccccccccccccccccccccccccc$	5	12.30	627798	delta-BHC	В	0.01885	197795.15		
7 13.55 623423 AldrinB 0.01985 195415.49 8 14.33 6036 B 0.00604 1772.36 9 14.99 623587 Hept. epoxideB 0.02017 188127.59 10 15.28 613362 gamma chlordaneB 0.01947 185797.48 11 15.58 576703 alpha chlordaneB 0.01922 178359.33 12 15.78 1167258 $4.4'$ -DDEB 0.03842 372407.92 13 17.43 979809 Endosulfan IIB 0.00528 1353.81 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.03468 173807.90 16 18.35 570396 Endrin aldebydeB	6	12.42	13472		V	0.01347	2967.81		
8 14.33 6036 B 0.00604 1772.36 9 14.99 623587 Hept. epoxide B 0.02017 188127.59 10 15.28 613362 gamma chlordane B 0.01947 185797.48 11 15.58 576703 alpha chlordane B 0.01947 185797.48 11 15.58 576703 alpha chlordane B 0.01922 178359.33 12 15.78 1167258 4,4'-DDE B 0.03842 372407.92 13 17.43 979809 Endosulfan II B 0.00528 1353.81 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.03468 193807.90 -12.69 16 18 35 570396 Endrin aldebyde B 0.03468 193807.90 -12.69	7	13.55	623423	Aldrin	В	0.01985	195415.49		
9 14.99 623587 Hept. epoxide B 0.02017 188127.59 SEP 2 9 2009 10 15.28 613362 gamma chlordane B 0.01947 185797.48 11 15.58 576703 alpha chlordane B 0.01922 178359.33 12 15.78 1167258 4,4'-DDE B 0.03842 372407.92 13 17.43 979809 Endosulfan II B 0.03948 272904.51 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.00528 1353.81 15 18.09 39898 B 0.00528 1353.81 0.00528 1353.81 0.003468 193807.90 -12 69.	8	14.33	6036		В	0.00604	1772.36		
10 15.28 613362 gamma chlordane B 0.01947 185797.48 4 11 15.58 576703 alpha chlordane B 0.01947 185797.48 4 11 15.58 576703 alpha chlordane B 0.01922 178359.33 4 12 15.78 1167258 $4.4'$ -DDE B 0.03942 372407.92 13 17.43 979809 Endosulfan II B 0.03948 272904.51 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.03468 193807.90 -12669 16 18 35 570396 Endrin aldebyde B 0.03468 193807.90 -12669	9	14.99	623587	Hept. epoxide	В	0.02017	188127.59	\sim	SEP 2 9 2009
11 15.58 576703 alpha chlordane B 0.01922 178359.33 12 15.78 1167258 $4,4'-DDE$ B 0.03842 372407.92 13 17.43 979809 Endosulfan II B 0.03948 272904.51 14 17.60 5283 B 0.00528 1353.81 15 18.09 39298 B 0.03999 10246.90 16 18.35 570396 Endrin aldebyde B 0.03468 173807.90 -12.69	10	15.28	613362	gamma chlordane	В	0.01947	185797.48	C.	5. Yan Ya Qin
12 15.78 1167258 4,4'-DDE B 0.03842 372407.92 13 17.43 979809 Endosulfan II B 0.03948 272904.51 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.039468 10246.90 16 18.35 570396 Endrin aldebyde B 0.03468 173807.90	11	15.58	576703	alpha chlordane	В	0.01922	178359.33	Compared and the second	
13 17.43 979809 Endosulfan II B 0.03948 272904.51 14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.039468 10246.90 16 18.35 570396 Endrin aldebyde B 0.03468 173807.90 -12.69	12	15.78	1167258	4,4'-DDE	В	0.03842	372407.92		
14 17.60 5283 B 0.00528 1353.81 15 18.09 39898 B 0.03990 10246.90 -1269	13	17.43	979809	Endosulfan II	₿	0.03948	272904.51		
15 18.09 39898 B $(0.03990 10246.90 -1269)$	14	17.60	5283		В	0.00528	1353.81		
16 18 35 576396 Endrin aldebyde B ($0.03468 183807.90 - 12 b L$)	15	18.09	39898		∖₽	0.03990	10246.90	(- car	
	16	18.35	576396	Endrin aldehyde	jз	(0.03468	133807.90	-12.66	

09/28/2009 15:39:15 Result: H:\TURBO6\6890-06\6-SEQ46\6A46221.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.		Height [µV]
17	19.30	670080	Endo. Sulfate	В	0.03711	181048.87
18	19.90	879709	Endrin ketone	В	0.03716	228958.87
19	20.27	3654		B	0.00365	1116.16
20	21.65	3778		В	0.00378	1119.46
21	22.09	936660		В	0.93666	223672.93
		9114664			1.81144	2.64e+06

TotalChrom Method File H:\TURBO6\6890-06\6B-(09-28-09).mth Printed by : NearyM on: 09/28/2009 14:39:23 Created by : NearyM on: 09/28/2009 14:28:33 Edited by : NearyM on: 09/28/2009 14:39:20 Number of Times Edited : 1 Number of Times Calibrated : 2641 Description: CURVE 07-23-09	Processed by: UM 9 128 109
Global Sample Information Default Sample Volume 1.000 ul Quantitation Units ng Void Time 0.000 min Correct amounts during calibration Yes Convert unknowns to concentration units Yes Reject outliers during calibration No An External Standard calibration will be used Units and the count tist of units and tist.	Reviewed by: SEP <u>F.9.2009</u>
Component Information alpha-BHC (2C) Component Type : Single Peak Component Retention Time : 12.289 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% User Values Label : Value 1 : 0.020000	0
Value 2 : 0.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Area Height ISTD Amt. ISTD Resp. # F	Replicates
A 0.0050 59167.60 18230.78 B 0.0100 120055.20 37653.06 C 0.0500 726062.00 230910.57 D 0.1000 1513058.90 479053.22 E 0.1500 2306418.00 729252.79	1 1 1 1 1
Calibration Curve : y = (-33889,937727) + (15537178.444673)x + (0.00000 R-squared : 0.999820 gamma-BHC (20) Component Type : Single Peak Component Retention Time : 13.247 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%	00)x^2 + (0.000000)x^3

Second sec

User Value Label Value 1 Value 2 Value 3 Value 3 Value 4 Value 5	s 0.020000 0.000000 0.000000 0.000000 0.000000							
Calibration Level Name	Level e Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	5	
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	59417.60 118588.20 685272.10 1409625.00 2120290.10	17741.79 35646.39 209785.75 431218.74 645746.54			1 1 1 1 1		
Calibration	Curve : v	- (-20106.8 3 999950	35296) + (14	1265800.0684	445)x + (0.00	0000)x^2 + ((0.000000)x^3	
beta-BHC (2C) Component Type Retention Time Search Window Search Window Efference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%								
User Value Label Value 1 Value 2 Value 3 Value 4 Value 5	s 0.020000 0.000000 0.000000 0.000000 0.000000	·						
Calibration Level Name	Level e Amount	Area	Height	ISTD Amt.	STD Resp. a	#Replicates		
A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	31441.00 60227.40 297272.80 582006.40 857706.90	9354.00 17632.68 88241.61 174733.28 259086.75			1 1 1 1		
Calibration R-squared	Curve : y : 0.	- (5894.68 5 999801	773) + (571	1685.738143)x + (0.00000)0)x^2 + (0.00	00000)x^3	
R-squared : 0.999801 delta-BHC (2C) Component Type : Single Peak Component Retention Time : 14.214 min Search Window : 3.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%								

	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.020000 0.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 1 0.1500 2	55688.70 113668.32 675666.54 1404673.29 2140504.20	16441.42 33795.96 203589.82 423766.22 643470.70			1 1 1 1	
	Calibration C R-squared		- (-29682.3 8 999834	8272) + (14)	408294.9834	412)x + (0.00	0000)x^2 + (0).000000)x^3
(H	leptachlor (2C Component T Retention Tir Search Wind Reference C Find peak clo Calibrating A Curve will igr Amounts will Weighting far Component s	c) Fype ne ow omponent osest to ex rea versus nore the or not be sca ctor for the standard p	: Single Pe : 14.398 m : 3.00 s, 0. : pected RT is s Amount us igin aled prior to o regression: urity percent	eak Compon in 00 % window ing a 1st On the regressi 1 tage : 100.0	der Fit on 000%			
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.020000 0.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 1 0.1500 1	61457.80 120302.80 654283.00 1325228.80 1983283.70	18066.65 35423.45 197797.65 399527.60 599015.06			1 1 1 1 1	
	Calibration C R-squared	Curve : y=	- (8613.04 3 999979	868) + (132	94035.46903	36)x + (0.000	000)x^2 + (0.	000000)x^3
	Idrin (2C) Component T Retention Tir Search Wind Reference C Find peak clo Calibrating A Curve will igr Amounts will Weighting fac Component s	Type ne owponent osest to ex rea versus nore the or not be sca ctor for the standard p	: Single Pe : 15.200 m : 5.00 s, 0. : pected RT in Amount us igin aled prior to regression: urity percent	eak Compon in 00 % n window ing a 1st Ord the regressi 1 tage : 100.0	ent der Fit on 000%			

	User Values Label : Value 1 : 0 Value 2 : 0 Value 3 : 0 Value 4 : 0 Value 5 : 0	.020000 .000000 .000000 .000000						
	Calibration Le	vel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	; -
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	54557.40 107511.80 602751.32 231225.16 844681.05	15912.56 31439.00 180007.00 366550.85 550737.14			1 1 1 1	
	Calibration Cu R-squared	$y = \frac{1}{2}$	(-12528.)6 99965	9154) + (12	391646.145	510)x + (0.00	0000)x^2 + ((0.000000)x^3
Ċ	lept. epoxide (2 Component Ty Retention Tim Search Windo Reference Co Find peak clos Calibrating Arc Curve will ign Amounts will r Weighting fact Component st	2C) ype w mponent sest to exp ea versus ore the origination to be scal tor for the landard pu	: Single Pe : 16.551 m : 5.00 s, 0. : bected RT in Amount us gin led prior to regression: inity percent	eak Compon in 00 % ing a 1st On the regressi 1 tage : 100.0	der Fit on 000%			
	User Values Label : Value 1 : 0 Value 2 : 5 Value 3 : 0 Value 4 : 0 Value 5 : 0	.500000 .000000 .000000 .000000						
	Calibration Le	vel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	56885.20 109701.80 570593.60 139066.60 701251.60	15948.20 30656.28 163889.29 327226.71 489068.98			1 1 1 1 1	
	Calibration Cu R-squared	arve : y =	(-354.6428 99980	01) + (1136	2768.298429	9)x + (0.0000	00)x^2 + (0.0	00000)x^3
9	amma chlordar Component Ty Retention Tim Search Windo Reference Co Find peak clos Calibrating Are Curve will igno Amounts will r Weighting fact Component st	ne (2C) ype mponent sest to exp ea versus ore the originate scal tor for the iandard put	: Single Pe : 16.989 m : 5.00 s, 0. : Dected RT in Amount us gin led prior to regression: rity percent	eak Compon in 00 % ing a 1st Or the regressi 1 tage : 100.0	ent der Fit on 000%			

	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000 0.000000						
	Calibration Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	54276.80 105280.70 564286.10 1136915.00 1702346.30	15444.82 30070.63 162750.42 329360.78 497516.11			1 1 1 1	
	Calibration C R-squared	urve y	= (-5225 ,808 999986	933) + (113	94393.4751	31)x + (0.000	000)x^2 + (0.0	00000)x^3
	Ipha chlordan Component 1 Retention Tir Search Wind Reference Ca Find peak clo Calibrating A Curve will igr Amounts will Weighting fac Component s User Values Label : Value 1 : Value 2 : Value 2 :	e (2O) Fype ow omponent osest to e: rea versu nore the o not be sc ctor for the standard p 0.040000 0.000000 0.000000	: Single Pe : 17.322 m : 3.00 s, 0 t : xpected RT i s Amount us rigin aled prior to e regression purity percen	eak Compon in 00 % ing a 1st Or the regressi 1 tage : 100.0	der Fit on 000%			
	Value 5 :	0.000000						
	Calibration Lo	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	53786.56 103223.44 535929.84 1069716.70 1602854.83	15001.26 29010.23 152244.87 307266.50 464401.83			1 1 1 1	
	Calibration C R-squared	Curve : y : 0.	= (-891.9267 999991	07) + (1069	8320.26184	5)x + (0.0000	00)x^2 + (0.00)0000)x^3
E	ndosulfan I (2 Component 1 Retention Tir Search Wind Reference Co Find peak clo Calibrating A Curve will igr Amounts will Weighting fac Component s	PC) Type ne ow omponent osest to ex rea versu nore the o not be sc ctor for the standard p	: Single Pe : 17.458 m : 5.00 s, 0. kpected RT i s Amount us rigin aled prior to e regression: purity percen	eak Compon in 00 % window ing a 1st On the regressi 1 tage : 100.0	der Fit on 000%			

	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	i
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	50821.44 97919.16 510027.66 1016266.30 1517438.37	14014.15 27015.10 142663.58 286681.09 428166.16			1 1 1 1	
	Calibration C R-squared	Curve : y	= (37 11687) 999976	3) + (101342	245.548093):	x + (0.00000	0)x^2 + (0.000	0000)x^3
4	4'-DDE (2C) Component Retention Ti Search Wind Reference C Find peak cl Calibrating A Curve will ig Amounts will Weighting fa Component	Type me dow component osest to ex Area versu nore the o I not be sc ictor for the standard p	: Single Pe : 17.672 m : 3.00 s, 0. : : : : : : : : : : : : :	eak Compon in 00 % ing a <u>1st Or</u> the regressi 1 tage : 100.0	der Fit ion 000%			
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	48053.80 94266.80 523132.60 1067699.60 1620306.00	14292.09 28282.82 159879.86 328370.68 497341.65			1 1 1 1 1	
	Calibration C R-squared	Curve : y : 0.	= (-13268.57 999922	75327) + (10	856513.259	162)x + (0.00	0000)x^2 + ((0.000000)x^3
	ieldrin (2C) Component Retention Ti Search Wind Reference C Find peak cl Calibrating A Curve will ig Amounts will Weighting fa Component	Type me dow component osest to ex Area versu nore the o I not be sc lotor for the standard p	: Single Pe : 18.075 m : 5.00 s, 0. : : s Amount us rigin aled prior to e regression: purity percent	eak Compon in 00 % ing a 1st Or the regressi 1 tage : 100.0	der Fit ion 000%			

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	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.040000 0.000000 0.000000 0.000000 0.000000							
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A B C D E	0.0050 0.0100 0.0500 0.1000 1 0.1500 1	51412.30 99979.00 545745.40 106901.00 672510.40	14029.96 27416.16 151976.87 309857.59 469130.73			1 1 1 1		
	Calibration C R-squared	Calibration Curve : y = (-10093.517111) + (11196874.783726)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999965							
Ċ	Endrin (2C) Component Type : Single Peak Component Retention Time : 18.757 min Search Window : 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%								
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.040000 0.000000 0.000000 0.000000 0.000000							
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates		
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500 1	41044.80 81892.80 451106.40 918414.20 400785.90	10945.91 21477.60 120598.52 248307.45 377473.76			1 1 1 1 1		
	Calibration C R-squared	Curve : y = : 0.9	(-12035)55 999879	2426) + (93	75942.07738	34)x + (0.000	000)x^2 + (0.	000000)x^3	
4	,4'-DDD (2C) Component Retention Til Search Wind Reference C Find peak cle Calibrating A Curve will ign Amounts will Weighting fa Component	Type me low component cosest to ex vrea versus nore the ori not be sca ctor for the standard pu	: Single Pe : 18.947 m : 5.00 s, 0. : pected RT in Amount us igin iled prior to regression: urity percent	eak Compon in 00 % ing a 1st Or the regressi 1 tage : 100.0	der Fit on 000%				

	User Values Label : Value 1 : 0 Value 2 : 0 Value 3 : 0 Value 4 : 0 Value 5 : 0	.040000 .000000 .000000 .000000 .000000						
	Calibration Le	vel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500 1	37298.60 73741.60 410304.80 834868.00 263363.40	10751.53 21005.91 119212.70 246217.27 374036.93			1 1 1 1 1	
	Calibration Cu R-squared	irve : y = : 0.9	<u>= (-95</u> 93.899 999956	856) + (846	8399.37128	7)x + (0.0000	00)x^2 + (0.0	00000)x^3
	ndosulfan II (2) Component Ty Retention Tim Search Windo Reference Con Find peak clos Calibrating Are Curve will igno Amounts will n Weighting fact Component st	e w mponent sest to ex sa versus ore the ori tot be sca tor for the andard p	: Single Pe : 19.216 m : 5.00 s, 0. : pected RT li Amount us igin aled prior to regression: urity percent	eak Compor in 00 % ing a 1st Or the regressi 1 tage : 100.0	der Fit on 000%			
	User Values Label : Value 1 : 0. Value 2 : 5. Value 3 : 0. Value 4 : 0. Value 5 : 0.	.500000 .000000 .000000 .000000 .000000						
	Calibration Lev Level Name	vel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500 1	45469.20 87730.00 448220.90 894340.40 340207.00	11894.82 22839.14 119255.84 240781.29 358781.10			1 1 1 1 1	
	Calibration Cu R-squared	rve : y =	(90.476734 999995	l) + (893814 ,	l3.226440)x	+ (0.000000)	x^2 + (0.0000	100)x^3
4	A DDT (2C) Somponent Ty Retention Time Search Windo Reference Cor Find peak clos Calibrating Are Curve will igno Amounts will n Weighting fact Component sta	ype e wmponent sest to exp ea versus or the ori iot be sca or for the andard pu	: Single Pe : 19.643 m : 5.00 s, 0. : pected RT in Amount us igin led prior to regression: urity percent	eak Compon in 00 % ing a 1st On the regressi 1 age : 100.0	der Eit on 000%			

	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.040000 0.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	31589.00 64793.20 365127.50 759959.40 1187841.42	9069.85 18358.57 105163.40 221005.90 345297.72			1 1 1 1	
	Calibration C B-equared	HINE : V	€ (-1887).8 ⁻ 999195	10944) + (79	48157.2352	20)x + (0.000	000)x^2 + (0.0)00000)x^3
	ndrin aldehyd Component Retention Tii Search Wind Reference C Find peak clu Calibrating A Curve will ig Amounts will Weighting fa Component	de (2C) Type me dow componen osest to e vrea versu nore the o nore the o nore the sc ctor for th standard p): Single P : 19.946 n : 5.00 s, 0 t : xpected RT(s Amount us rigin aled prior to e regression purity percer	eak Compor nin .00 % in window sing a 1st Or the regressi : 1 ntage : 100.0	der Fit ion 000%			
	User Values Label : Value 1 : Value 2 : Value 3 : Value 4 : Value 5 :	0.500000 5.000000 0.000000 0.000000 0.000000						
	Calibration L Level Name	evel Amount	Area	Height	ISTD Amt.	ISTD Resp. #	# Replicates	
	A B C D E	0.0050 0.0100 0.0500 0.1000 0.1500	33689.00 63994.80 309576.02 605698.30 900022.38	8619.24 - 16146.27 - 78644.33 - 156988.57 - 235535.57 -			1 1 1 1 1	
	Calibration C R-squared	Curve : y : 0.	= (5987.172 999926	926) + (5977	7919.456170))x + (0.00000	00)x^2 + (0.00	0000)x^3
E	ndo. Sulfate Component Retention Til Search Wind Reference C Find peak cle Calibrating A Curve will ig Amounts will Weighting fa Component	(20) Type low component ossest to ex rea versu nore the o not be so ctor for th standard p	: Single P : 20.556 n : 5.00 s, 0 t : xpected RT s Amount us rigin aled prior to e regression burity percen	eak Compor nin .00 % sing a 1st Or the regressi : 1 itage : 100.0	der Fit			

User Values Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area A 0.0050 29240.70 7655.72 --В 61952.00 15862.73 -----0.0100 С 0.0500 337930.20 87416.66 --1 D 0.1000 690907.62 181730.62 ------1 Ε 0.1500 1078513.00 284891.28 -----1 Calibration Curve : y = (-13914.812750) + (7200373.280367)x + (0.000000)x^2 + (0.000000)x^3 : 0.999347 R-squared Methoxychlor (2C) Component Type Single Peak Component Retention Time 21.085 min Search Window 5.00 s, 0.00 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000% **User Values** Label Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000 Calibration Level Level Name Amount Height ISTD Amt. ISTD Resp. # Replicates Area A 0.0050 17846.80 5063.86 -----____ 1 В 36919.20 10243.22 -----0.0100 1 С 0.0500 186268.70 51370.70 ------1 0.1000 368121.10 101902.68 ------D 1 Ε 0.1500 556530.00 155237.14 ------1 Calibration Curve $\cdot y = (-330.885340) + (3705841.989529)x + (0.000000)x^2 + (0.000000)x^3$ R-squared 0.999965 Endrin ketone (2C) Component Type Single Peak Component Retention Time 21.777 min 5.00 s, 0.00 % Search Window Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
Ă	0.0050	41785.90	9413.88			1
В	0.0100	82589.40	18305.33			1
С	0.0500	427450.40	95982.14			1
D	0.1000	858780.30	196672.77			1
E	0.1500	1337743.20	305060.08			1

Calibration Curve : $y = (-10242.057624) + (8887490.438482)x + (0.000000)x^2 + (0.000000)x^3$ R-squared (0.999397)

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98759	Date	: 09/28/2009 14:35:22
Operator	: tchrom	Sample Name	: 9090862-CCV1
Sample Number	: 0.15	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/15
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 11:30:02	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46215.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46215.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46215.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46215.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46215.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
2	8.69	1535		В	0.00154	501.06
3	9.99	11264		В	0.01126	3314.13
4	10.95	1583	•	В	0.00158	527.69
5	12.30	2306418	alpha-BHC (2C)	В	0.15000	729252.79
7	13.26	2120290	gamma-BHČ (2C)	V	0.15000	645746.54
8	13.48	857707	beta-BHC (2C)	В	0.15000	259086.75
9	14.04	13868		В	0.01387	3797.14
10	14.23	2140504	delta-BHC (2C)	В	0.15000	643470.70
11	14.41	1983284	Heptachlor (2C)	В	0.15000	599015.06
12	14.65	4460		В	0.00446	1216.54
13	15.21	1844681	Aldrin (2C)	В	0.15000	550737.14
14	15.35	1663		V	0.00166	605.93
16	16.15	15384		В	0.01538	4491.06
17	16.56	1701252	Hept. epoxide (2C)	В	0.15000	489068.98
18	17.00	1702346	gamma chlordane (2C)	В	0.15000	497516.11
19	17.34	1602855	alpha chlordane (2C)	В	0.15000	464401.83

09/28/2009 14:35:22 Result: H:\TURBO6\6890-06\6-SEQ46\6b46215.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
20	17.47	1517438	Endosulfan I (2C)	v	0.15000	428166.16
21	17.69	1620306	4,4'-DDE (2C)	В	0.15000	497341.65
22	18.09	1672510	Dieldrin (2C)	В	0.15000	469130.73
24	18.77	1400786	Endrin (2C)	В	0.15000	377473.76
25	18.96	1263363	4,4'-DDD (2C)	В	0.15000	374036.93
26	19.23	1340207	Endosulfan II (2C)	В	0.15000	358781.10
27	19.47	5084	, ,	В	0.00508	845.02
.28	19.65	1187841	4,4'-DDT (2C)	V	0.15000	345297.72
29	19.96	900022	Endrin aldehyde (2C)	В	0.15000	235535.57
30	20.19	43955		V	0.04395	8157.03
31	20.57	1078513	Endo. Sulfate (2C)	В	0.15000	284891.28
32	20.76	2887		В	0.00289	703.68
33	21.09	556530	Methoxychlor (2C)	В	0.15000	155237.14
34	21.37	4005		В	0.00401	1098.50
35	21.56	7417		В	0.00742	1763.63
36	21.79	1337743	Endrin ketone (2C)	В	0.15000	305060.08
37	23.51	1701		В	0.00170	135.40
		30249403			3.11481	8.74e+06

Software Version	: 6.2.1.0.104:0104	Date	: 09/28/2009 14:35:29
Reprocess Number	: but2048: 98761		
Operator	: tchrom	Sample Name	: 9090863-CCV1
Sample Number	: 0.10	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/16
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 12:05:55	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46216.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46216.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46216.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46216.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46216.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.28	1887		В	0.00189	722.67
2	8.67	2094		В	0.00209	691.32
3	9.98	16055		В	0.01606	4727.28
4	10.94	2796		В	0.00280	915.41
5	12.29	1513059	alpha-BHC (2C)	В	0.10000	479053.22
7	13.25	1409625	gamma-BHC (2C)	В	0.10000	431218.74
8	13.46	582006	beta-BHC (2C)	В	0.10000	174733.28
9	14.03	10862		В	0.01086	2899.26
10	14.21	1404673	delta-BHC (2C)	V	0.10000	423766.22
11	14.40	1325229	Heptachlor (2Ć)	В	0.10000	399527.60
12	14.64	2772	,	В	0.00277	780.29
13	15.20	1231225	Aldrin (2C)	В	0.10000	366550.85
14	15.34	1745		V	0.00175	604.24
15	16.14	10013		В	0.01001	2973.90
16	16.55	1139067	Hept. epoxide (2C)	В	0.10000	327226.71
17	16.99	1136915	gamma chlordane (2C)	В	0.10000	329360.78

09/28/2009 14:35:29 Result: H:\TURBO6\6890-06\6-SEQ46\6b46216.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	17.32	1069717	alpha chlordane (2C)	В	0.10000	307266.50
19	17.46	1016266	Endosulfan I (2C)	V	0.10000	286681.09
20	17.67	1067700	4,4'-DDE (2C)	В	0.10000	328370.68
21	18.08	1106901	Dieldrin (2C)	В	0.10000	309857.59
22	18.76	918414	Endrin (2C)	В	0.10000	248307.45
23	18.95	834868	4,4'-DDD (2C)	В	0.10000	246217.27
24	19.22	894340	Endosulfan II (2C)	В	0.10000	240781.29
.25,	19.64	759959	4,4'-DDT (2C)	В	0.10000	221005.90
26	19.95	605698	Endrin aldehyde (2C)	В	0.10000	156988.57
27	20.17	61678		V	0.06168	12860.43
28	20.56	690908	Endo. Sulfate (2C)	В	0.10000	181730.62
29	20.74	1845		V	0.00185	474.56
30	21.08	368121	Methoxychlor (2C)	В	0.10000	101902.68
31	21.36	2922		В	0.00292	815.93
32	21.55	5393		В	0.00539	1276.17
33	21.78	858780	Endrin ketone (2C)	В	0.10000	196672.77
		20053535			2.12006	5.79e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98763	Date	: 09/28/2009 14:35:36
Operator	: tchrom	Sample Name	: 9091289-CCV1
Sample Number	: 0.05	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/17
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	e : 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.00000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 12:41:53	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46217.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46217.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46217.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46217.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46217.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak	Time	Area	Component	BL		Height
#	funul	[uv-sec]	Name		CONCENTRATION	Įμνj
1	8.28	1538		В	0.00154	624.83
3	9.98	10248		В	0.01025	2948.82
4	10.94	1741		В	0.00174	572.50
5	12.29	726062	alpha-BHC (2C)	В	0.05000	230910.57
6	13.25	685272	gamma-BHC (2C)	В	0.05000	209785.75
7	13.46	297273	beta-BHC (2C)	В	0.05000	88241.61
8	14.03	6622		В	0.00662	1696.59
9	14.21	675667	delta-BHC (2C)	V	0.05000	203589.82
10	14.40	654283	Heptachlor (2C)	В	0.05000	197797.65
11	15.20	602751	Aldrin (2C)	В	0.05000	180007.00
12	15.34	1714		V	0.00171	600.03
13	16.14	4872		В	0.00487	1441.53
14	16.55	570594	Hept. epoxide (2C)	В	0.05000	163889.29
15	16.99	564286	gamma chlordane (2C)	В	0.05000	162750.42
16	17.32	535930	alpha chlordane (2C)	В	0.05000	152244.87
17	17.46	510028	Endosulfan I (2C)	V	0.05000	142663.58

09/28/2009 14:35:36 Result: H:\TURBO6\6890-06\6-SEQ46\6b46217.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	17.67	523133	4,4'-DDE (2C)	В	0.05000	159879.86
19	18.08	545745	Dieldrin (2C)	В	0.05000	151976.87
20	18.76	451106	Endrin (2C)	В	0.05000	120598.52
21	18.95	410305	4,4'-DDD (2C)	В	0.05000	119212.70
22	19.22	448221	Endosulfan II (2C)	В	0.05000	119255.84
23	19.64	365128	4.4'-DDT (2C)	В	0.05000	105163.40
24	19.95	309576	Endrin aldehyde (2C)	В	0.05000	78644.33
25	20.17	35768		V	0.03577	7773.43
26	20.56	337930	Endo. Sulfate (2C)	В	0.05000	87416.66
27	21.09	186269	Methoxychlor (2C)	В	0.05000	51370.70
28	21.55	2143		В	0.00214	562.70
29	21.78	427450	Endrin ketone (2C)	В	0.05000	95982.14
		9891653			1.06465	2.84e+06

Software Version	: 6.2.1.0.104:0104	Date	: 09/28/2009 14:35:42
Reprocess Number	: buf2048: 98765		
Operator	: tchrom	Sample Name	: 9091538-CCV1
Sample Number	: 0.01	Study	: PEST ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/18
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.00000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 13:17:47	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46218.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46218.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46218.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46218.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46218.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	0.00	2075			0.00208	007.92
2	12.33	120055	alpha, BHC (2C)	u p	0.002.90	27652.06
2	12.20	120000		D D	0.01000	37033.00
3	13.20	110000	gamma-Bric (2C)	B	0.01000	30040.39
4	13.47	60227	beta-BHC (2C)	В	0.01000	17632.68
5	14.04	1850		В	0.00185	490.68
6	14.22	113668	delta-BHC (2C)	V	0.01000	33795.96
7	14.40	120303	Heptachlor (2C)	В	0.01000	35423.45
8	15.20	107512	Aldrin (2C)	В	0.01000	31439.00
9	16.56	109702	Hept. epoxide (2C)	В	0.01000	30656.28
10	17.00	105281	gamma chlordane (2C)	В	0.01000	30070.63
11	17.33	103223	alpha chlordane (2C)	В	0.01000	29010.23
12	17.46	97919	Endosulfan I (2C)	V	0.01000	27015.10
13	17.68	94267	4,4'-DDE (2C)	В	0.01000	28282.82
14	18.08	99979	Dieldrin (2C)	В	0.01000	27416.16
15	18.76	81893	Endrin (2C)	В	0.01000	21477.60
16	18.95	73742	4.4'-DDD (2C)	В	0.01000	21005.91

09/28/2009 14:35:42 Result: H:\TURBO6\6890-06\6-SEQ46\6b46218.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
17	19.22	87730	Endosulfan II (2C)	в	0.01000	22839.14
18	19.65	64793	4,4'-DDT (2C)	В	0.01000	18358.57
19	19.95	63995	Endrin aldehyde (2C)	В	0.01000	16146.27
20	20.18	8428		В	0.00843	1986.12
21	20.56	61952	Endo. Sulfate (2C)	В	0.01000	15862.73
22	21.09	36919	Methoxychlor (2C)	В	0.01000	10243.22
23	21.78	82589	Endrin ketone (2C)	В	0.01000	18305.33
		1817591			0.21325	511665.17

Software Version	: 6.2.1.0.104:0104 : buf2048: 98767	Date	: 09/28/2009 14:35:48
Operator Sample Number AutoSampler Instrument Name Instrument Serial #	: tchrom : 0.005 : BUILT-IN : HP6890-06 : CN10520010	Sample Name Study Rack/Vial Channel A/D mV Range	: 9091537-CCV1 : PEST ICAL : 1/19 : B : 1000 : 20.50 min
Sampling Rate Sample Volume Sample Amount Data Acquisition Time	 0.00 min 5.0000 pts/s 1.000000 ul 1.0000 09/28/2009 13:53:55 	Area Reject Dilution Factor Cycle	: 29.50 min : 1500.000000 : 1.00 : 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46219.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46219.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46219.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46219.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46219.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	9.99	1681		В	0.00168	564.35
2	12.30	59168	alpha-BHC (2C)	В	0.00500	18230.78
3	13.25	59418	gamma-BHČ (2C)	В	0.00500	17741.79
4	13.47	31441	beta-BHC (2C)	В	0.00500	9354.00
5	14.22	55689	delta-BHC (2C)	В	0.00500	16441.42
6	14.40	61458	Heptachlor (2C)	В	0.00500	18066.65
7	15.21	54557	Aldrin (2C)	В	0.00500	15912.56
8	16.56	56885	Hept. epoxide (2C)	В	0.00500	15948.20
9	17.00	54277	gamma chlordane (2C)	В	0.00500	15444.82
10	17.33	53787	alpha chlordane (2C)	В	0.00500	15001.26
11	17.47	50821	Endosulfan I (2C)	V	0.00500	14014.15
12	17.68	48054	4,4'-DDE (2C)	В	0.00500	14292.09
13	18.08	51412	Dieldrin (2C)	В	0.00500	14029.96
14	18.77	41045	Endrin (2C)	В	0.00500	10945.91
15	18.96	37299	4,4'-DDD (2C)	В	0.00500	10751.53
16	19.22	45469	Endosulfan II (2C)	В	0.00500	11894.82

09/28/2009 14:35:48 Result: H:\TURBO6\6890-06\6-SEQ46\6b46219.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	19.65	31589	4,4'-DDT (2C)	В	0.00500	9069.85
18	19.95	33689	Endrin aldehyde (2C)	В	0.00500	8619.24
19	20.18	4928		В	0.00493	1227.66
20	20.56	29241	Endo. Sulfate (2C)	В	0.00500	7655.72
21	21.09	17847	Methoxychlor (2C)	В	0.00500	5063.86
22	21.78	41786	Endrin ketone (2C)	В	0.00500	9413.88
		921539			0.10661	259684.50

Software Version	: 6.2.1.0.104:0104	Date	: 09/28/2009 14:59:56
Reprocess Number	: DUI2046: 96769	0	0004500
Operator	: tchrom	Sample Name	: 9091536
Sample Number	: 0.05	Study	: ALT PEST
AutoSampler	: BUILT-IN	Rack/Vial	: 1/20
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 14:29:42	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46220.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46220.rst

Inst Method : h:\turbo6\6890-06\6890-66 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46220.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46220.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46220.rst Report Format File: h:\turbo6\6890-06\06\06\07.01

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.29	BB	775151	alpha-BHC (2C)	0.05207	247698.29	4.1	12.24 -	12.34
13.25	BB	730576	gamma-BHC (2C)	0.05262	222961.47	5.2	13.20 -	13.30
13.47	BB	315628	beta-BHC (2C)	0.05423	93583.58	8.5	13.42 -	13.52
14.22	VB	721298	delta-BHC (2C)	0.05212	218695.34	4.2	14.17 -	14.27
14.40	BB	698323	Heptachlor (2C)	0.05318	210138.63	6.4	14.35 -	14.45
15.20	ΒV	658835	Aldrin (2C)	0.05418	195387.61	8.4	15.15 -	15.25
16.55	BB	612069	Hept. epoxide (2C)	0.05390	175208.17	7.8	16.50 -	16.60
16.99	BB	589482	gamma chlordane (2C)	0.05219	170091.59	4.4	16.94 -	17.04
17.33	ΒV	571505	alpha chlordane (2C)	0.05350	162931.03	7.0	17.28 -	17.38
17.46	VB	540382	Endosulfan I (2C)	0.05332	151728.11	6.6	17.41 -	17.51
17.68	BB	556831	4,4'-DDE (2C)	0.05251	168740.63	5.0	17.63 -	17.73
18.08	BΒ	571728	Dieldrin (2C)	0.05196	159471.55	3.9	18.03 -	18.13
18.76	BB	471577	Endrin (2C)	0.05158	127953.29	3.2	18.71 -	18.81
18.95	BB	440232	4,4'-DDD (2C)	0.05312	128451.96	6.2	18.90 -	19.00
19.22	BΒ	478225	Endosulfan II (2C)	0.05349	127188.19	7.0	19.17 -	19.27
19.65	BB	406381	4,4'-DDT (2C)	0.05350	117293.35	7.0	19.60 -	19.70

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09/28/2009 14:59:56 Result: H:\TURBO6\6890-06\6-SEQ46\6b46220.rst

Ret Time [min]	BL.	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window	- Relative
19.95 20.56 21.09 21.78	BV BB BB BB	396768 355975 200517 455895	Endrin aldehyde (2C) Endo. Sulfate (2C) Methoxychlor (2C) Endrin ketone (2C)	0.06537 0.05137 0.05420 0.05245	102332.27 92278.04 54742.12 102679.19	30.7 2.7 8.4 4.9	 19.90 -20.51 21.04 21.73 	- 20.00 - 20.61 - 21.14 - 21.83
10547378 Missing Component Report Component Expected Retention (Calibration File) All components were found					3.03e+06		e (w) xc a b	etyde)

SEP 29 2009

Software Version Reprocess Number	: 6.2.1.0.104:0104 · buf2048: 98783	Date	: 09/28/2009 15:39:18
Operator	: tchrom	Sample Name	: 9040971
Sample Number		Study	: INDBM MIX
AutoSampler	: BUILT-IN	Rack/Vial	: 1/21
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 15:09:19	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46221.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46221.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46221.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46221.rst Calib Method : h:\turbo6\6890-06\6b-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46221.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]	
1	8.28	2154		в	0.00215	789.88	() ~ ~ ~ ~
2	9.99	4560		В	0.00456	1290.79	(Se or o and a and
3	10.53	186087		В	0.18609	57791.62	SEL 7 A TOUR
4	13.47	115953	beta-BHC (2C)	в	0.01927	34197.17	Carren
5	14.04	5565		В	0.00556	1346.38	
6	14.22	246214	delta-BHC (2C)	V	0.01915	73679.02	GA LOLAND
7	15.20	233202	Aldrin (2C)	В	0.01983	68875.58	N WWW
8	15.34	2374		V	0.00237	754.56	A
9	16.15	1913		В	0.00191	572.16	qh
10	16.56	230366	Hept. epoxide (2C)	в	0.02030	64609.97	
11	16.99	221472	gamma chlordane (2C)	В	0.01990	62750.53	
12	17.33	213183	alpha chlordane (2C)	В	0.02001	59870.93	
13	17.68	407707	4,4'-DDE (2C)	В	0.03878	123716.95	
14	19.22	353241	Endosulfan II (2C)	В	_0.03951	93100.90	6.00
15	19.95	217415	Endrin aldehyde (2C)	В	0.03537	∕55325.34	F10,5 %)
16	20.18	22104		V	0.02210	4696.22	

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09/28/2009 15:39:18 Result: H:\TURBO6\6890-06\6-SEQ46\6b46221.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	20.56	260224	Endo. Sulfate (2C)	В	0.03807	67122.65
18	21.78	328353	Endrin ketone (2C)	В	0.03810	73973.56
19	23.54	2405		В	0.00241	155.40
20	25.22	367062		В	0.36706	62019.88
		3421553			0.90251	906639.51

TotalChrom Method File H:\TU	RBO6\6890-06\6a-SURR-(09-28-09).mth
Printed by	: NearyM on: 09/29/2009 07:41:51
Created by	: NearyM on: 09/29/2009 07:32:39
Edited by	: NearyM on: 09/29/2009 07:41:46
Number of Times Edited	: 1
Number of Times Calibrated	: 2656
Description: CURVE 07-23-0)9
Global Sample Informatio	n
Default Sample Volume	: 1.000 ul
Quantitation Units	: ng
Void Time	: 0.000 min

Quantitation Units		ng	
Void Time	:	0.000 min	
Correct amounts during calibration	:	Yes	
Convert unknowns to concentration units	:	Yes	
Reject outliers during calibration	:	No	

An External Standard calibration will be used

Unknown peaks will be quantitated using a response factor of 1.000000e+06 First peak will be relative retention reference

Component Information

Tetrachloro-m-xylene Component Type : Single Peak Component Retention Time : 9.178 min Search Window : 7.00 s, 0.50 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	119169.00	37708.67			1
В	0.0100	232827.30	73832.26			1
С	0.0500	1250293.00	404388.28	***********		1
D	0.0750	1927388.00	622331.90			1
E	0.1000	2577555.80	837450.45			1

Calibration Curve : $y = (-23852.108321) + (25943723.506686)x + (0.000000)x^2 + (0.000000)x^3$ R-squared 0.999825

Decachlorobiphenyl

Component Type : Single Peak Component Retention Time 22.082 min Search Window 7.00 s, 0.50 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1

Component standard purity percentage : 100.0000%



09/29/2009 07:41:51 Method: H:\TURBO6\6890-06\6a-SURR-(09-28-09).mth

User Values Label : Value 1 : 0,500000 Value 2 : 5,000000 Value 3 : 0,000000 Value 4 : 0,000000 Value 5 : 0,000000

Calibration Level Level Name Amount Area Height ISTD Amt. ISTD Resp. # Replicates Â 0.0050 128712.70 31295.74 -1 0.0100 242343.40 58421.16 -----В 1 С 0.0500 1139436.40 277998.06 ------1 D E 0.0750 1714333.60 416324.66 1 0.1000 2250104.70 557451.05 ------1

Calibration Curve : $y = (19246.754829) + (22411237.607727)x + (0.000000)x^2 + (0.000000)x^3 R-squared : 0.999906$

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98788	Date	: 09/29/2009 07:38:55
Operator	: tchrom	Sample Name	: 9090851
Sample Number	:	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/22
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.00000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 15:45:08	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46222.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46222.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46222.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46222.rst Calib Method : h:\turbo6\6890-06\6a-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46222.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	4024		В	0.00402	1539.26
2	9.18	2577556	Tetrachloro-m-xylene	В	0.10000	837450.45
3	10.28	27548		В	0.02755	8694.39
4	10.41	4273		V	0.00427	1343.60
5	12.04	12675		В	0.01267	3964.44
6	15.27	3611		В	0.00361	1175.29
7	15.74	10109		В	0.01011	940.98
8	17.60	3484		В	0.00348	764.44
9	21.65	4503		В	0.00450	1224.55
10	22.08	2250105	Decachlorobiphenyl	В	0.10000	557451.05
		4897887			0.27023	1.41e+06
Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98790	Date	: 09/29/2009 07:39:02			
--------------------------------------	--	-----------------	-----------------------			
Operator	: tchrom	Sample Name	: 9090852			
Sample Number	:	Study	: SURR ICAL			
AutoSampler	: BUILT-IN	Rack/Vial	: 1/23			
Instrument Name	: HP6890-06	Channel	: A			
Instrument Serial #	: CN10520010	A/D mV Range	: 1000			
Delay Time	: 0.00 min	End Time	: 29.50 min			
Sampling Rate	: 5.0000 pts/s					
Sample Volume	: 1.000000 ul	Area Reject	: 1500.00000			
Sample Amount	: 1.0000	Dilution Factor	: 1.00			
Data Acquisition Time	: 09/28/2009 16:21:11	Cycle	: 2			

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46223.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46223.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46223.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46223.rst Calib Method : h:\turbo6\6890-06\6a-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46223.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	3265		В	0.00327	1257.79
2	9.18	1927388	Tetrachloro-m-xylene	В	0.07500	622331.90
3	10.27	20811	-	В	0.02081	6632.07
4	12.04	9698		В	0.00970	3019.31
5	15.88	16465		В	0.01647	332.43
6	17.73	8462		В	0.00846	290.59
7	22.08	1714334	Decachlorobiphenyl	В	0.07500	416324.66
		3700423			0.20870	1.05e+06

Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:39:08
Reprocess Number	: buf2048: 98792		
Operator	: tchrom	Sample Name	: 9090254
Sample Number	:	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/24
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 16:57:02	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46224.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46224.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46224.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46224.rst Calib Method : h:\turbo6\6890-06\6a-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46224.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.75	4019		В	0.00402	1552.76
2	9.17	1250293	Tetrachloro-m-xylene	В	0.05000	404388.28
3	10.27	13834	·	В	0.01383	4384.93
4	12.03	6207		В	0.00621	1999.92
5	15.26	4751		В	0.00475	1545.13
6	15.78	9458		В	0.00946	486.72
7	17.71	8157		В	0.00816	378.03
8	22.08	1139436	Decachlorobiphenyl	B	0.05000	277998.06
		2436156			0.14643	692733.81

Software Version Reprocess Number	: 6.2.1.0.104:0104 / buf2048: 98794	Date	: 09/29/2009 07:39:14
Operator	: tchrom	Sample Name	: 9091567
Sample Number	•	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/25
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 17:33:00	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46225.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46225.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46225.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46225.rst Calib Method : h:\turbo6\6890-06\6a-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46225.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	9.18	232827	Tetrachloro-m-xylene	в	0.01000	73832.26
2	13.31	1874	-	В	0.00187	441.36
3	14.96	1603		В	0.00160	421.21
4	15.26	5873		В	0.00587	1906.90
5	15.90	16303		В	0.01630	5453.39
7	16.95	69072		В	0.06907	23372.53
9	22.08	242343	Decachlorobiphenyl	В	0.01000	58421.16
		569896			0.11473	163848.82

Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:39:19
Reprocess Number	: but2048: 98796		0001500
Operator	: tchrom	Sample Name	: 9091568
Sample Number	:	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/26
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 18:09:00	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46226.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46226.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46226.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46226.rst Calib Method : h:\turbo6\6890-06\6a-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46226.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	9.18	119169	Tetrachloro-m-xylene	В	0.00500	37708.67
3	13.30	1909	-	В	0.00191	563.43
4	14.98	5695		В	0.00569	1227.17
5	15.26	9276		В	0.00928	2879.47
6	15.90	20090		В	0.02009	6500.46
7	16.43	2728		В	0.00273	1013.87
8	16.95	82411		В	0.08241	27412.17
11	22.08	128713	Decachlorobiphenyl	В	0.00500	31295.74
		369990			0.13211	108600.99

FotalChrom Method File H:\TU	R	BO6\689(D-061	6B-SURR-(0	9-28-09).mth
Printed by	:	NearyM	on:	09/29/2009	07:42:32
Created by	:	NearyM	on:	09/29/2009	07:33:44
Edited by	:	NearyM	on:	09/29/2009	07:42:27
Number of Times Edited	:	1			
Number of Times Calibrated	:	2641			
Description: CURVE 07-23-0	90				

Global Sample Information

Default Sample Volume	:1.000 ul
Quantitation Units	: ng
Void Time	: 0.000 min
Correct amounts during calibration	: Yes
Convert unknowns to concentration units	: Yes
Reject outliers during calibration	: No

An External Standard calibration will be used

Unknown peaks will be quantitated using a response factor of 1.000000e+06 First peak will be relative retention reference

Component Information

Tetrachloro-m-xylene (2C) Component Type : Single Peak Component Retention Time : 10.530 min Search Window : 7.00 s, 0.50 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1 Component standard purity percentage : 100.0000%

User Values

Label : Value 1 : 0.500000 Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates
A	0.0050	46383.70	14157.08			1
В	0.0100	88928.00	26974.68			1
С	0.0500	463292.40	144422.53		····	1
D	0.0750	714487.80	225268.81			1
E	0.1000	952166.50	300154.47			1

Calibration Curve : $y = (-5955.632689) + (9562652.347697)x + (0.000000)x^{2} + (0.000000)x^{3}$ R-squared : 0.999815

Decachlorobiphenyl (2C) Component Type : Single Peak Component Retention Time : 25.216 min Search Window : 7.00 s, 0.50 % Reference Component : Find peak closest to expected RT in window Calibrating Area versus Amount using a 1st Order Fit Curve will ignore the origin Amounts will not be scaled prior to the regression Weighting factor for the regression: 1

Component standard purity percentage : 100.0000%

Processed by

09/29/2009 07:42:32 Method: H:\TURBO6\6890-06\6B-SURR-(09-28-09).mth

User Values Label : Value 1 : 0.500000 Value 2 : 5.000000

Value 2 : 5.000000 Value 3 : 0.000000 Value 4 : 0.000000 Value 5 : 0.000000

Calibration Level

Level Name	Amount	Area	Height	ISTD Amt.	ISTD Resp.	# Replicates	
				· · · · · · · · · · · · · · · · · · ·	···		
А	0.0050	55059.20	9412.32		<u> </u>	1	
В	0.0100	101855.50	17168.59			1	
С	0.0500	448912.60	75947.00			1	
D	0.0750	660225.60	111823.85	***********	************	1	
E	0.1000	855718.50	145686.58			1	

Calibration Curve $y = (17936.041902) + (8467046.627043)x + (0.000000)x^2 + (0.000000)x^3$ R-squared (: 0.999548

Software Version Reprocess Number	: 6.2.1.0.104:0104 · buf2048· 98789	Date	: 09/29/2009 07:38:59
Operator	: tchrom	Sample Name	: 9090851
Sample Number	•	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/22
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 15:45:08	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46222.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46222.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46222.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46222.rst Calib Method : h:\turbo6\6890-06\6b-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46222.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	2555		В	0.00256	958.06
2	10.53	952167	Tetrachloro-m-xylene (В	0.10000	300154.47
3	11.62	10373	2	В	0.01037	3188.12
5	13.76	5210		В	0.00521	1543.38
6	15.35	2222		В	0.00222	721.27
8	25.22	855719	Decachlorobiphenyl (2C	В	0.10000	145686.58
		1828244			0.22036	452251.87

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98791	Date	: 09/29/2009 07:39:05
Operator	: tchrom	Sample Name	: 9090852
Sample Number		Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/23
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 16:21:11	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46223.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46223.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46223.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46223.rst Calib Method : h:\turbo6\6890-06\6b-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46223.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	8.28	2032		В	0.00203	782.37
2	10.53	714488	Tetrachloro-m-xylene (В	0.07500	225268.81
3	11.61	7703	2 .	В	0.00770	2333.89
4	13.76	4012		В	0.00401	1191.52
5	15.34	1966		В	0.00197	644.24
7	23.52	2725		В	0.00273	200.73
8	25.21	660226	Decachlorobiphenyl (2C	В	0.07500	111823.85
		1393151			0.16844	342245.41

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98793	Date	: 09/29/2009 07:39:11
Operator	; tchrom	Sample Name	: 9090254
Sample Number		Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/24
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 16:57:02	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46224.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46224.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46224.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46224.rst Calib Method : h:\turbo6\6890-06\6b-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46224.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL.		Height [µV]
1	8.28	2614		В	0.00261	959.16
2	10.53	463292	Tetrachloro-m-xylene (В	0.05000	144422.53
3	11.61	4974	•	В	0.00497	1549.23
4	13.75	2502		В	0.00250	769.22
5	15.34	2844		В	0.00284	889.46
7	25.21	448913	Decachlorobiphenyl (2C	В	0.05000	75947.00
		925139			0.11293	224536.60

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98795	Date	: 09/29/2009 07:39:17
Operator	: tchrom	Sample Name	: 9091567
Sample Number		Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/25
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 17:33:00	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46225.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46225.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46225.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46225.rst Calib Method : h:\turbo6\6890-06\6b-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46225.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	10.53	88928	Tetrachloro-m-xvlene (В	0.01000	26974.68
2	15.35	3129		В	0.00313	966.02
3	16.77	5309		В	0.00531	1849.56
4	17.34	26340		В	0.02634	8830.67
5	25.22	101856	Decachlorobiphenyl (2C	₿	0.01000	17168.59
		225561			0.05478	55789.53

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98797	Date	: 09/29/2009 07:39:22
Operator	: tchrom	Sample Name	· 9091568
Sample Number	:	Study	: SURR ICAL
AutoSampler	: BUILT-IN	Rack/Vial	: 1/26
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 18:09:00	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46226.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46226.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46226.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46226.rst Calib Method : h:\turbo6\6890-06\6b-surr-(09-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46226.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq

300 Response [mV] 200 SPL1 40.0 100 ____ 10.53 -25.22 -17.34 16.76 15.34 19.91 TETRAC DECAC Harrison Har 12 ż à 6 8 10 14 16 18 20 22 24 26 28 Time [min]

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	10.53	46384	Tetrachloro-m-xylene (В	0.00500	14157.08
2	15.34	3325		В	0.00333	1051.26
3	16.76	6319		В	0.00632	2197.69
4	17.34	31464		В	0.03146	10498.29
6	25.22	55059	Decachlorobiphenyl (2C	В	0.00500	9412.32
		142551			0.05111	37316.64

9/25/09 Nm

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98508	Date	: 09/25/2009 08:17:24
Operator	: tchrom	Sample Name	: 9090864-CCV2
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/76
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 01:59:14	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46176.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46176.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46176.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46176.rst

Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46176.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
10.94	BB	1809285	alpha-BHC	0.06075	593528.82	21.5	10.89 -	10.99
11.79	VB	1634067	gamma-BHC	0.05991	518398.24	19.8	11.74 -	11.84
12.03	BB	678526	beta-BHC	0.06102	210064.87	22.0	11.98 -	12.08
12.48	BB	1660958	delta-BHC	0.05958	520664.23	19.2	12.43 -	12.53
13.00	BB	1602279	Heptachlor	0.06249	502022.19	25.0	12.95 -	13.05
13.73	BB	1526683	Aldrin	0.06236	475848.51	24.7	13.68 -	13.78
15.18	BB	1477954	Hept. epoxide	0.05952	444138.34	19.0	15.13 -	15.23
15.47	BB	1512498	gamma chlordane	0.05741	459590.23	14.8	15.42 -	15.52
15.77	BB	1434712	alpha chlordane	0.05928	428845.05	18.6	15.72 -	15.82
15.97	ΒV	1415663	4,4'-DDE	0.05550	442803.85	11.0	15.92 -	16.02
16.08	VB	1331935	Endosulfan I	0.05594	387382.45	11.9	16.03 -	16.13
16.62	BB	1457390	Dieldrin	0.05335	429237.57	6.7	16.57 -	16.67
17.13	BΒ	1269438	Endrin	0.05198	366144.71	4.0	17.08 -	17.18
17.29	BB	1061333	4,4'-DDD	0.04893	327357.41	-2.1	17.24 -	17.34
17.62	BB	1159125	Endosulfan II	0.05023	324150.73	0.5	17.57 -	17.67
17.90	BB	964102	4,4'-DDT	0.04287	290786.26	-14.3	17.85 -	17.95
18.54	BB	764171	Endrin aldehyde	0.04413	206923.19	-11.7	18.49 -	18.59

09/25/2009 08:17:24 Result: H:\TURBO6\6890-06\6-SEQ46\6A46176.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
18.94	VB	478991	Methoxychlor	0.04095	141374.29	-18.1	18.89 -	18.99
19.49	BB	979151	Endo. Sulfate	0.04629	264498.42	-7.4	19.44 -	19.54
20.10	BB	1119695	Endrin ketone	0.04570	292901.31	-8.6	20.05 -	20.15
22.58	ВB	30951	Decachlorobiphenyl	7.64e-04	260.06	- 98.5 -	22.53 -	22.63
		25368909		1.07895	7.63e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene

9.380

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 08:17:31
Reprocess Number	: but2048: 98510		
Operator	: tchrom	Sample Name	: 9090254-CCV2
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/77
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	9 : 09/25/2009 02:35:12	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46177.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46177.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46177.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46177.rst Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46177.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
9.36	BB	1174746	Tetrachloro-m-xylene	0.06974	379597.70	39.5	9.31 -	9.41
15.45	BΒ	8273	ganhma chlordane	0.00213	2494.50	-9 57 .7	15.40 -	15.50
15.96	BΒ	6282	4, 4-D DE	0.00263	310.50	-94.7	15.91 -	16.01
17.88	BΒ	7316	4,4'-DDT	0.00243	377.32	-95.1	17.83 -	17.93
22.32	BB	1082008	Decachlorobiphenyl	0.04769	252357.13	-4.6	22.27 -	22.37
				0.40400				
		22/8625		0.12463	635137.15			

9/25/09 mm

Missing Component Report

Component Expected Retention (Calibration File)

alpha-BHC	10.938
gamma-BHC	11.787
beta-BHC	12.034
delta-BHC	12.478
Heptachlor	12.999
Aldrin	13.731
Hept. epoxide	15.180
alpha chlordane	15.770
Endosulfan I	16.075

9/25/09

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98526	Date	: 09/25/2009 08:18:19
Operator	: tchrom	Sample Name	: 9090864-CCV3
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/85
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 07:22:47	Cycle	: 10

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46185.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46185.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46185.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46185.rst

Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46185.rst Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
10.93	BB	1828997	alpha-BHC	0.06137	602288.22	22.7	10.88 -	10.98
11.78	VB	1654229	gamma-BHC	0.06061	528556.31	21.2	11.73 -	11.83
12.03	BB	681911	beta-BHC	0.06132	210868.48	22.6	11.98 -	12.08
12.47	BB	1662254	delta-BHC	0.05963	519299.47	19.3	12.42 -	12.52
12.99	BB	1591899	Heptachlor	0.06210	502122.44	24.2	12.94 -	13.04
13.72	BB	1531300	Aldrin	0.06254	476415.12	25.1	13.67 -	13.77
15.17	BB	1481805	Hept. epoxide	0.05967	446986.13	19.3	15.12 -	15.22
15.45	BB	1511702	gamma chlordane	0.05738	463435.12	14.8	15.40 -	15.50
15.76	BB	1400820	alpha chlordane	0.05792	422881.85	15.8	15.71 -	15.81
15.96	ΒV	1423056	4,4'-DDE	0.05577	450692.37	11.5	15.91 -	16.01
16.07	VB	1332866	Endosulfan I	0.05598	390442.64	12.0	16.02 -	16.12
16.61	BB	1456404	Dieldrin	0.05331	435000.96	6.6	16.56 -	16.66
17.12	BB	1235485	Endrin	0.05062	355713.64	1.2	17.07 -	17.17
17.28	BB	1050704	4,4'-DDD	0.04846	320822.41	-3.1	17.23 -	17.33
17.62	BB	1165663	Endosulfan II	0.05051	326326.34	1.0	17.57 -	17.67
17.89	BB	934590	4,4'-DDT	0.04162	281198.70	-16.8	17.84 -	17.94
18.54	BB	778689	Endrin aldehyde	0.04496	210916.40	-10.1	18.49 -	18.59

09/25/2009 08:18:19 Result: H:\TURBO6\6890-06\6-SEQ46\6A46185.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
18.93	VB	456573	Methoxychlor	0.03903	136075.58	-21.9	18.88 -	18.98
19.48	BB	986418	Endo. Sulfate	0.04663	269006.63	-6.7	19.43 -	19.53
20.09	BB	1121291	Endrin ketone	0.04576	295262.77	-8.5	20.04 -	20.14
22.27	BB	13859	Decachlorobiphenyl	8.72e-07	1313.37	-100.0	22.22 -	22.32
		25300515		1.07520	7.65e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene

9.380

a/25/09

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 10:25:13
Reprocess Number	: buf2048: 98573		
Operator	: tchrom	Sample Name	: 9090254-CCV3
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/86
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	e : 09/25/2009 07:58:22	Cycle	; 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46186.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46186.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46186.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46186.rst

Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46186.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
9.35	BB	1202664	Tetrachloro-m-xylene	0.07135	386112.95	42.7	9.30 -	9.40
15.45	BB	5124	gamma chlordane	0.00201	1593.34	-96,0	15.40 -	15.50
15.96	BB	5968	4,4'-DDE	0.00262	322.82	-94.8	15.91 -	16.01
17.88	ΒB	8658		0.00249	386.09	-95.0	17.83 -	17.93
22.31	BB	1103068	Decachlorobiphenyl	0.04863	256742.04	-2.7	22.26 -	22.36
		2325481		0.12711	645157.25			

Missing Component Report

Component Expected Retention (Calibration File)

alpha-BHC	10.938
gamma-BHC	11.787
beta-BHC	12.034
delta-BHC	12.478
Heptachlor	12.999
Aldrin	13.731
Hept. epoxide	15.180
alpha chlordane	15.770
Endosulfan I	16.075

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 10:25:18
Reprocess Number	: buf2048: 98575		
Operator	: tchrom	Sample Name	: 9090636
Sample Number	:	Study	: PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/87
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 08:34:23	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46187.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46187.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46187.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46187.rst

Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46187.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL.	NG CONCENTRATION	Height [µV]
1	9.35	496430	Tetrachloro-m-xylene	В	0.03050	161137.94
2	10.93	307394	alpha-BHC	В	0.01320	100320.79
3	11.78	299746	gamma-BHC	В	0.01348	94515.74
4	12.03	128444	beta-BHC	В	0.01249	38594.91
5	15.96	7978	4,4'-DDE	В	0.00269	2457.74
6	16.71	2989		В	0.00299	999.09
7	16.90	17138		В	0.01714	5199.68
8	17.12	1230746	Endrin	В	0.05043	353957.01
9	17.29	23875	4,4'-DDD	В	0.00291	6354.74
10	17.89	2067122	4,4'-DDT	В	0.08948	631070.64
11	18.43	5424		В	0.00542	1732.99
12	18.54	30957	Endrin aldehyde	V	0.00220	8466.92
13	18.80	29587	•	В	0.02959	6940.35

9/25/09 pum 5.370 DDT 1.570 DDT 1.570

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09/25/2009 10:25:18 Result: H:\TURBO6\6890-06\6-SEQ46\6A46187.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
<u></u>				·		
14	18.93	2566191	Methoxychior	V	0.21929	772680.19
15	19.11	7772	•	V	0.00777	2446.14
16	20.09	37418	Endrin ketone	В	0.00268	9699.53
17	22.31	474456	Decachlorobiphenyl	В	0.02057	111463.46
		7733667			0.52282	2.31e+06

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 14:56:14
Reprocess Number	: but2048: 98618		
Operator	: tchrom	Sample Name	: 9090864-CCV4
Sample Number	;	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/96
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	e : 09/25/2009 14:07:21	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46196.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46196.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46196.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46196.rst Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46196.rst Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
10.94	BB	1706345	alpha-BHC	0.05749	560640.22	15.0	10.89 -	10.99
11.79	VB	1567566	gamma-BHC	0.05760	499860.21	15.2	11.74 -	11.84
12.04	BB	650323	beta-BHC	0.05854	201331.22	17.1	11.99 -	12.09
12.48	BB	1596797	delta-BHC	0.05742	504195.94	14.8	12.43 -	12.53
13.00	BB	1549260	Heptachlor	0.06048	487898.65	21.0	12.95 -	13.05
13.73	BB	1468662	Aldrin	0.06008	462023.74	20.2	13.68 -	13.78
15.18	BB	1431724	Hept. epoxide	0.05770	429028.89	15.4	15.13 -	15.23
15.47	BB	1465085	gamma chlordane	0.05567	448470.49	11.3	15.42 -	15.52
15.77	BB	1300242	alpha chlordane	0.05387	393173.95	7.7	15.72 -	15.82
15.97	ΒV	1381873	4,4'-DDE	0.05423	440312.27	8.5	15.92 -	16.02
16.08	VB	1289687	Endosulfan I	0.05420	376414.10	8.4	16.03 -	16.13
16.62	BB	1416885	Dieldrin	0.05192	417991.95	3.8	16.57 -	16.67
17.13	BB	1201872	Endrin	0.04927	346610.34	-1.5	17.08 -	17.18
17.29	BB	1033807	4,4'-DDD	0.04771	317980.69	-4.6	17.24 -	17.34
17.63	BB	1134420	Endosulfan II	0.04918	318464.86	-1.6	17.58 -	17.68
17.90	BΒ	948156	4,4'-DDT	0.04219	282854.14	-15.6	17.85 -	17.95
18.55	BB	751160	Endrin aldehyde	0.04338	201722.67	-13.2	18.50 -	18.60

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09/25/2009 14:56:14 Result: H:\TURBO6\6890-06\6-SEQ46\6A46196.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
18.94	VB	466360	Methoxychlor	0.03987	138848.53	-20.3	18.89 -	18.99
19.49	BB	956843	Endo. Sulfate	0.04526	256907.93	-9.5	19.44 -	19.54
20.10	BB	1101068	Endrin ketone	0.04496	287621.34	-10.1	20.05 -	20.15
22.31	BB	15957	Decachlorobiphenyl	9.45e-05	753.30	-99.8	22.26 -	22.36
		24434093		1.04111	7.37e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene

9.380

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 15:34:51
Reprocess Number	: buf2048: 98630		
Operator	: tchrom	Sample Name	: 9090254-CCV4
Sample Number		Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/97
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 14:53:10	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46197.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46197.rst

Result File : H:\10RB06\6890-06\6890-06\6890-6 12-5-08-ins from H:\TURB06\6890-06\6-SEQ46\6A46197.raw Inst Method : h:\turbo6\6890-06\6890-06\6a-process.mth from H:\TURB06\6890-06\6-SEQ46\6A46197.rst Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURB06\6890-06\6-SEQ46\6A46197.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	ΒL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
			· · · · · · · · · · · · · · · · · · ·					
9.36	BΒ	1131951	Tetrachloro-m-xylene	0.06726	367053.31	(34.5)	9.31 -	9.41
15.45	BB	7179	gamma chlordane	0.00209	2223.94	-95.8	15.40 -	15.50
15.93	BB	6663	4,4'-DDE	0.00265	554.13	-94.7	15.88 -	15.98
17.89	BB	11375	4,4'-DDT	0.00261	515.16	94.8	17.84 -	17.94
22.31	BB	1112097	Decachlorobiphenyl	0.04904	258588.43	(-1.9)	22.26 -	22.36
				- <u></u>		\mathcal{C}		
		2269265		0.12364	628934.96			

Missing Component Report

Component	Expected Retention	(Calibration File)
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alpha-BHC	10.938
gamma-BHC	11.787
beta-BHC	12.034
delta-BHC	12.478
Heptachlor	12.999
Aldrin	13.731
Hept. epoxide	15.180
alpha chlordane	15.770
Endosulfan I	16.075

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98509	Date	: 09/25/2009 08:17:28
Operator	: tchrom	Sample Name	: 9090864-CCV2
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/76
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 01:59:14	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46176.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46176.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46176.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46176.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46176.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.48	BB	684791	alpha-BHC (2C)	0.04796	215878.06	-4.1	12.43 -	12.53
13.44	VB	636191	gamma-BHC (2C)	0.04967	193038.90	-0.7	13.39 -	13.49
13.66	BB	278715	beta-BHC (2C)	0.05163	82224.96	3.3	13.61 -	13.71
14.41	BB	643941	delta-BHC (2C)	0.05144	192237.50	2.9	14.36 -	14.46
14.60	BΒ	611817	Heptachlor (2C)	0.05310	184580.48	6.2	14.55 -	14.65
15.40	ΒV	571253	Aldrin (2C)	0.05468	169727.59	9.4	15.35 -	15.45
16.76	BΒ	529549	Hept. epoxide (2C)	0.05531	149972.23	10.6	16.71 -	16.81
17.20	BB	505277	gamma chlordane (2C)	0.05327	145534.51	6.5	17.15 -	17.25
17.53	ΒV	495587	alpha chlordane (2C)	0.05470	140732.08	9.4	17.48 -	17.58
17.67	VB	462919	Endosulfan I (2C)	0.05462	127956.44	9.2	17.62 -	17.72
17.87	BB	472528	4,4'-DDE (2C)	0.05414	143284.15	8.3	17.82 -	17.92
18.29	BB	495553	Dieldrin (2C)	0.05409	137630.24	8.2	18.24 -	18.34
18.97	BB	413678	Endrin (2C)	0.05580	110582.76	11.6	18.92 -	19 <i>.</i> 02
19.15	BΒ	375701	4,4'-DDD (2C)	0.05344	108128.58	6.9	19.10 -	19.20
19.43	BB	411575	Endosulfan II (2C)	0.05443	108593.98	8.9	19.38 -	19.48
19.85	BB	285317	4,4'-DDT (2C)	0.04506	82252.18	-9.9	19.80 -	19.90

9/25/09 12/10/

¹⁹⁹/225 e 2 of 2

09/25/2009 08:17:28 Result: H:\TURBO6\6890-06\6-SEQ46\6b46176.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - F	Relative
20.16 20.77 21.29 22.02	BB BB BB BB	280370 365302 148598 388407	Endrin aldehyde (2C) Endo. Sulfate (2C) Methoxychlor (2C) Endrin ketone (2C)	0.05040 0.05166 0.04368 0.04803	71900.57 93877.48 39343.52 85146.92	0.8 3.3 -12.6 -3.9	20.11 - 20.72 - 21.24 - 21.97 -	20.21 20.82 21.34 22.07
		9057071		1.03710	2.58e+06			

Missing Component Pepor	ŧ
Component	Expected Retention (Calibration File)
component	

Tetrachloro-m-xylene (2C)	10.720
Decachlorobiphenyl (2C)	25.700

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98511	Date	: 09/25/2009 08:17:35
Operator	: tchrom	Sample Name	: 9090254-CCV2
Sample Number		Study	: SURR
AutoSampler	; BUILT-IN	Rack/Vial	: 1/77
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 02:35:12	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46177.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46177.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46177.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46177.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46177.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - Re	elative	
10.71	BB	422344	Tetrachloro-m-xylene (0.04394	132067.64	-12.1	10.66 -	10.76	9[1
20.00	00	838978		0.10060	200573.71	13.5	23.30 -	20.00	Ç -

Missing Component Report

Component Expected Detention (Colibration E			
Component Expected Retention (Calibration F	File)	e)	1

alpha-BHC (2C)	12.480
gamma-BHC (2C)	13.440
beta-BHC (2C)	13.657
delta-BHC (2C)	14.410
Heptachlor (2C)	14.597
Aldrin (2C)	15.403
Hept. epoxide (2C)	16.756
gamma chlordane (2C)	17.195
alpha chlordane (2C)	17.529
Endosulfan I (2C)	17.667
4,4'-DDE (2C)	17.874
Dieldrin (2C)	18.286

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 08:18:21
Reprocess Number	: buf2048: 98527		
Operator	: tchrom	Sample Name	: 9090864-CCV3
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/85
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 07:22:47	Cycle	: 10

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46185.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46185.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46185.raw

Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46185.rst

Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46185.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.47	BB	695674	alpha-BHC (2C)	0.04867	220460.05	-2.7	12.42 -	12.52
13.43	VB	649780	gamma-BHC (2C)	0.05068	197853.18	1.4	13.38 -	13.48
13.65	BB	282109	beta-BHC (2C)	0.05226	83187.72	4.5	13.60 -	13.70
14.40	BB	654222	delta-BHC (2Ć)	0.05221	197142.60	4.4	14.35 -	14.45
14.59	BB	619874	Heptachlor (2C)	0.05378	186810.49	7.6	14.54 -	14.64
15.39	BV	580903	Aldrin (2C)	0.05557	173982.49	11.1	15.34 -	15.44
16.75	BB	539837	Hept, epoxide (2C)	0.05637	153740.14	12.7	16.70 -	16.80
17.19	BB	514066	gamma chlordane (2C)	0.05418	147984.60	8.4	17.14 -	17.24
17.52	ΒV	503587	alpha chlordane (2C)	0.05558	142915.85	11.2	17.47 -	17.57
17.66	VB	471085	Endosulfan I (2C)	0.05557	131162.59	11.1	17.61 -	17.71
17.87	BB	481529	4.4'-DDE (2C)	0.05513	145986.32	10.3	17.82 -	17.92
18.28	BB	504221	Dieldrin (2C)	0.05501	140344.24	10.0	18.23 -	18.33
18.96	BB	408748	Endrin (2C)	0.05515	109263.61	10.3	18.91 -	19.01
19.15	BB	382432	4.4'-DDD (2C)	0.05437	110315.06	8.7	19.10 -	19.20
19.42	BB	418279	Endosulfan II (2C)	0.05532	110416.82	10.6	19.37 -	19.47
19.84	BB	292206	4,4'-DDT (2C)	0.04610	83969.19	-7.8	19.79 -	19.89

9/25/09

09/25/2009 08:18:21 Result: H:\TURBO6\6890-06\6-SEQ46\6b46185.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
20.15 20.76 21.29 22.01	BB BB BB VB	287167 373213 149293 400297	Endrin aldehyde (2C) Endo. Sulfate (2C) Methoxychlor (2C) Endrin ketone (2C)	0.05164 0.05279 0.04389 0.04950	73058.90 95781.52 39755.45 87739.42	3.3 5.6 -12.2 -1.0	20.10 - 20.71 - 21.24 - 21.96 -	20.20 20.81 21.34 22.06
Missing C	omn	9208523		1.05377	2.63e+06			

missing component report	L
Component	Expected Retention (Calibration File)
\mathbf{T} the difference of \mathbf{r} (0.0)	40 700

Tetrachloro-m-xylene (2C)	10.720
Decachlorobiphenyl (2C)	25.700

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 10:25:16
Reprocess Number	: buf2048: 98574		
Operator	: tchrom	Sample Name	: 9090254-CCV3
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/86
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 07:58:22	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46186.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46186.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46186.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46186.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46186.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative	
10.71	BB	432459	Tetrachloro-m-xylene (0.04497	133908.32	-10.1	10.66 - 25.48 -	10.76	9psi
20.00	00	856752	Decaeniorosiphenyi (20	0.10270	203744.19	10.0	20.10	20.00	N

Missing Component Report

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Course o court	Even a start Data attar	(Calibratian File)	
Component	Expected Releniion	a subtation rules	
oomponone		(Ounoradon r no)	
		• • • • • • • • • • • • • • • • • • • •	

alpha-BHC (2C)	12.480
gamma-BHC (2C)	13.440
beta-BHC (2C)	13.657
delta-BHC (2C)	14.410
Heptachlor (2C)	14.597
Aldrin (2C)	15.403
Hept. epoxide (2C)	16.756
gamma chlordane (2C)	17.195
alpha chlordane (2C)	17.529
Endosulfan I (2C)	17.667
4,4'-DDE (2C)	17.874
Dieldrin (2C)	18.286

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98576	Date	: 09/25/2009 10:25:21
Operator	: tchrom	Sample Name	: 9090636
Sample Number		Study	: PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/87
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 08:34:23	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46187.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46187.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46187.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46187.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46187.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.44	1638		В	0.00164	606.76
2	10.71	180530	Tetrachloro-m-xylene (В	0.01950	55783.40
3	12.47	119834	alpha-BHC (2C)	В	0.01101	37504.25
4	13.43	121287	gamma-BHC (2C)	В	0.01141	36609.03
5	13.65	55100	beta-BHC (2C)	В	0.01005	16095.13
7	17.87	3171	4,4'-DDE (2C)	В	0.00228	919.82
8	18.96	410861	Endrin (2C)	В	0.05543	108797.22
9	19.15	21688	4,4'-DDD (2C)	V	0.00467	3948.38
10	19.84	638131	4,4'-DDT (2C)	В	0.09827	184306.93
11	20.15	14045	Endrin aldehyde (2C)	В	0.00184	3291.10
12	20.85	15024	Endo. Sulfate (2C)	В	0.00151	2959.04
13	20.95	10144		V	0.01014	2087.34

09/25/2009 10:25:21 Result: H:\TURBO6\6890-06\6-SEQ46\6b46187.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
14	21.29	812253	Methoxychlor (2C)	В	0.24096	220215.89
15	21.58	1712		В	0.00171	513.86
16	22.01	12765	Endrin ketone (2C)	В	0.00170	2966.19
17	25.54	192591	Decachlorobiphenyl (2C	В	0.02516	31684.83
		2610776			0.49728	708289.17

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98619	Date	: 09/25/2009 14:56:17
Operator Sample Number AutoSampler Instrument Name Instrument Serial # Delay Time Sampling Rate Sample Volume Sample Amount	t tchrom E BUILT-IN E HP6890-06 CN10520010 C0.00 min 5.0000 pts/s 1.000000 ul 1.0000	Sample Name Study Rack/Vial Channel A/D mV Range End Time Area Reject Dilution Factor	: 9090864-CCV4 : 8081 : 1/96 : B : 1000 : 29.48 min : 3000.000000 : 1.00
Data Acquisition Time	e : 09/25/2009 14:07:21	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46196.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46196.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46196.raw

Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46196.rst

Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46196.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.48	BB	670667	alpha-BHC (2C)	0.04704	212294.46	-5.9	12.43 -	12.53
13.44	VB	625249	gamma-BHC (2C)	0.04885	190764.34	-2.3	13.39 -	13.49
13.66	BB	272736	beta-BHC (2C)	0.05052	80533.10	1.0	13.61 -	13.71
14.41	BB	630729	delta-BHC (2C)	0.05044	190186.71	0.9	14.36 -	14.46
14.60	BB	605645	Heptachlor (2C)	0.05258	182878.85	5.2	14.55 -	14.65
15.40	ΒV	558119	Aldrin (2C)	0.05347	167148.63	6.9	15.35 -	15.45
16.76	ΒV	468241	Hept. epoxide (2C)	0.04900	133506.55	-2.0	16.71 -	16.81
17.20	VV	516414	gamma chlordane (2C)	0.05442	110186.66	8.8	17.15 -	17.25
17.53	٧V	439530	alpha chlordane (2C)	0.04851	101893.20	-3.0	17.48 -	17.58
17.67	VV	459625	Endosulfan I (2C)	0.05424	98873.70	8.5	17.62 -	17.72
17.88	VB	412039	4,4'-DDE (2C)	0.04745	112966.37	-5.1	17.83 -	17.93
18.29	BB	493378	Dieldrin (2C)	0.05386	136612.79	7.7	18.24 -	18.34
18.97	BB	398783	Endrin (2C)	0.05384	107044.07	7.7	18.92 -	19.02
19.15	BB	371863	4,4'-DDD (2C)	0.05291	107618.70	5.8	19.10 -	19.20
19.43	BΒ	405156	Endosulfan II (2C)	0.05358	108163.51	7.2	19.38 -	19.48
19.85	BΒ	284018	4,4'-DDT (2C)	0.04487	81759.07	-10.3	19.80 -	19.90

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09/25/2009 14:56:17 Result: H:\TURBO6\6890-06\6-SEQ46\6b46196.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
20.16 20.77 21.30 22.02	BB BB BB BB	271629 360928 155681 389991	Endrin aldehyde (2C) Endo. Sulfate (2C) Methoxychlor (2C) Endrin ketone (2C)	0.04881 0.05103 0.04579 0.04823	68844.05 93448.48 41350.56 85814.22	-2.4 2.1 -8.4 -3.5	20.11 - 20.72 - 21.25 - 21.97 -	20.21 20.82 21.35 22.07
		8790418		1.00943	2.41e+06			

Missing Component Report

Component Repor	t Expected Retention (Calibration File)
Tetrachloro-m-xylene (2C)	10.720
Decachlorobiphenyl (2C)	25.700

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 15:34:53
Reprocess Number	: buf2048: 98631		
Operator	: tchrom	Sample Name	: 9090254-CCV4
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/97
Instrument Name	; HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.44 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	; 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 14:53:10	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46197.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46197.rst Inst Method : h:\turbo6\6890-06\6890-66\2-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46197.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46197.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46197.rst

Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



	Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RI Window -	Relative
-	10.71 25.54	BB BB	436082 425305	Tetrachloro-m-xylene (Decachlorobiphenyl (2C	0.04533 0.05788	135574.24 70049.92	-9.3 15.8	10.66 - 25.49 -	10.76 25.59
			861387		0 10321	205624 16	\leq		

Missing Component Report

Component Expected Retention (Calibration File)

alpha-BHC (2C)	12.480
gamma-BHC (2C)	13.440
beta-BHC (2C)	13.657
delta-BHC (2C)	14.410
Heptachlor (2C)	14.597
Aldrin (2C)	15.403
Hept. epoxide (2C)	16.756
gamma chlordane (2C)	17.195
alpha chlordane (2C)	17.529
Endosulfan I (2C)	17.667
4,4'-DDE (2C)	17.874
Dieldrin (2C)	18.286

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9/29/09 3.070 3.070 5.070 5.070 5.070

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Software Version	: 6.2.1.0.104:0104 : buf2048: 98798	Date	: 09/29/2009 07:49:12
Operator	: tchrom	Sample Name	: 9091475-CCV1
Sample Number		Study	: PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/27
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 18:44:59	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46227.raw <Modified>



Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46227.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46227.rst

Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46227.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	3419		в	0.00342	1325.23
2	7.96	2262		В	0.00226	1033.66
3	8.67	3386		В	0.00339	1466.66
4	9.18	517077	Tetrachloro-m-xylene	В	0.02085	168259.52
5	9.37	4532	•	В	0.00453	1739.11
6	10.06	3440		В	0.00344	1424.93
7	10.75	333995	alpha-BHC	В	0.00952	109289.00
8	11.38	3394	•	В	0.00339	1420.89
9	11.60	325887	gamma-BHC	В	0.01020	103000.73
10	11.85	137869	beta-BHC	В	0.00915	41728.86
11	12.01	3266		В	0.00327	1372.53
12	12.62	3793		В	0.00379	1541.18
13	13.21	3760		В	0.00376	1574.99
14	13.78	4478		В	0.00448	1800.80
15	14.33	6463		В	0.00646	2350.45
16	14.87	5058		В	0.00506	2054.80

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09/29/2009 07:49:12 Result: H:\TURBO6\6890-06\6-SEQ46\6A46227.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.	NG CONCENTRATION	Height [µV]
17	15.27	3875	gamma chlordane	В	9.90e-04	1258.52
18	15.38	6512		V	0.00651	2139.43
19	15.78	36362	4,4'-DDE	в	0.00240	6606.21
20	15.88	7084	Endosulfan I	V	3.35e-04	2441.25
21	16.37	3873	Dieldrin	В	0.00106	1737.59
22	16.72	24786		в	0.02479	7430.95
23	16.93	1398509	Endrin	В	0.05211	398973.18
24	17.10	23588	4,4'-DDD	V	0.00205	6939.35
25	17.30	12533		V	0.01253	2769.07
26	17.49	4510	Endosulfan II	В	4.78e-04	1721.04
27	17.70	2558714	4,4'-DDT	В	0.10646	773776.08
28	18.18	3857		В	0.00386	1561.98
29	18.34	13500	Endrin aldehyde	В	0.00101	3973.78
30	18.61	15414		В	0.01541	4192.42
31	18.75	3071396	Methoxychlor	V	0.26672	926522.21
32	18.93	11783		V	0.01178	3684.98
33	19.01	3101		V	0.00310	1281.28
34	19.41	4145		В	0.00415	1565.07
35	19.80	2339		В	0.00234	1032.07
36	19.89	30064	Endrin ketone	V	0.00323	7916.63
37	20.18	1992		В	0.00199	938.19
38	20.55	3338		В	0.00334	1259.23
39	20.91	2633		В	0.00263	1065.25
40	21.27	2321		В	0.00232	949.38
41	21.63	5136		В	0.00514	1635.12
42	22.08	538738	Decachlorobiphenyl	В	0.02318	122890.75
		9152182			0.65688	2.73e+06

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Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98800	Date	: 09/29/2009 07:49:23
Operator	: tchrom	Sample Name	: 9090864-CCV1
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/28
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 19:21:06	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46228.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46228.rst

Inst Method : h:\turbo6\6890-06\6890-06\6890-06\6890-06\6890-06\6890-06\6-SEQ46\6A46228.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46228.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46228.rst Report Format File: h:\turbo6\6890-06\06%d.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time	BL	Area	Component	NG	Height	%D	RT Window -	Relative
[min]		[uV-sec]	Name	CONC.	[µV]	0.05ng		
10.75	BB	1925324	alpha-BHC	0.04990	636097.53	-0.2	10.70 -	10.80
11.60	BB	1757421	gamma-BHC	0.04991	560817.69	-0.2	11.55 -	11.65
11.85	BB	727021	beta-BHC	0.05075	225293.08	1.5	11.80 -	11.90
12.29	ΒE	1780126	delta-BHC	0.04990	559427.79	-0.2	12.24 -	12.34
12.81	BB	1710701	Heptachior	0.05134	540330.64	2.7	12.76 -	12.86
13.54	BB	1651351	Aldrin	0.05106	519104.34	2.1	13.49 -	13.59
14.98	VB	1592758	Hept. epoxide	0.05160	476163.38	3.2	14.93 -	15.03
15.27	BB	1639222	gamma chlordane	0.05058	501019.44	1.2	15.22 -	15.32
15.57	BB	1538268	alpha chlordane	0.05065	471595.95	1.3	15.52 -	15.62
15.77	ΒV	1542738	4,4'-DDE	0.05038	491821.93	0.8	15.72 -	15.82
15.87	VB	1425998	Endosulfan I	0.05116	420810.22	2.3	15.82 -	15.92
16.41	BB	1576035	Dieldrin	0.05067	466554.20	1.3	16.36 -	16.46
16.93	BB	1339529	Endrin	0.04997	385821.53	-0.1	16.88 -	16.98
17.09	BB	1182788	4,4'-DDD	0.05104	367398.53	2.1	17.04 -	17.14
17.42	BB	1263216	Endosulfan II	0.05081	352399.02	1.6	17.37 -	17,47
17.70	BB	1142363	4,4'-DDT	0.04897	349109.58	-2.1	17.65 -	17.75
18.34	BB	840619	Endrin aldehyde	0.05101	226701.73	2.0	18.29 -	18.39


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09/29/2009 07:49:23 Result: H:\TURBO6\6890-06\6-SEQ46\6A46228.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - F	Relative
18.75	BB	561833	Methoxychlor	0.04975	168498.59	-0.5	18.70 -	18.80
19.29	BB	924156	Endo. Sulfate	0.05010	252536.53	0.2	19.24 -	19.34
19.89	BB	1208524	Endrin ketone	0.05030	318171.36	0.6	19.84 -	19.94
22.08	BB	13079	Deeachlorobiphenyl	-2.8e-04	1048.99	-1 00.6 -	22.03 -	22.13
		27343069		1.00956	8.29e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene

9.178

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Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98802	Date	: 09/29/2009 07:49:33
Operator	: tchrom	Sample Name	: 9090254-CCV1
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/29
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.46 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 19:57:04	Cycle	: 3

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46229.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46229.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46229.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46229.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46229.rst Report Format File: h:\turbo6\6890-06\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative	
9.18 15.27	BB BB	1238009 4568	Tetrachloro-m-xylene	0.04864	399991.00 1503.46	-2.7 -98 2 0	9.13 - 15.22 -	9.23 15.32	9/29/09
15.89	BB	9484	Endosulfan I-	4.21e-04	276.46	-992	15.84 -	15.94	m
22.09	BB	1140969	Decachlorobiphenyl	0.05005	275175.15	0.1	22.04 -	22.14	//0.

0.10012 676946.07

Missing Component Report

2393031

C	Compone	nt E	Expected	Retention	(Calibration	File)

10.743
11.590
11.839
12.280
12.799
13.529
14.977
15.566
15.768
16.409

Page 1 of 2

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98804	Date	: 09/29/2009 07:49:43
Operator	: tchrom	Sample Name	: 9090308
Sample Number	:	Study	: TECH
AutoSampler	: BUILT-IN	Rack/Vial	: 1/30
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 20:32:59	Cycle	: 4

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46230.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46230.rst

Inst Method : h:\turbo6\6890-06\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46230.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46230.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46230.rst Report Format File: h:\turbo6\6890-06\6asmp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
1	7.28	3832		В	0.00383	1383.23
2	7.76	3065		В	0.00306	1229.72
3	8.99	166968		В	0.16697	54521.88
4	10.89	24584		В	0.02458	7985.45
5	11.17	19454		В	0.01945	6182.65
6	11.43	52482		В	0.05248	15401.29
7	11.74	12444		В	0.01244	4012.99
8	11.89	15288	beta-BHC	В	4.90e-04	4803.72
9	12.15	42184		В	0.04218	12980.08
10	12.25	33373	delta-BHC	V	0.00284	9691.34
11	12.40	82411		V	0.08241	20202.44
12	12.50	4624		V	0.00462	1626.97
13	12.61	442586		V	0.44259	137880.10
14	12.81	927086	Heptachlor	В	0.02793	292944.74
15	13.06	52932	-	V	0.05293	16015.99
16	13.12	135368		V	0.13537	23912.51

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09/29/2009 07:49:43 Result: H:\TURBO6\6890-06\6-SEQ46\6A46230.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
17	13.32	45717		V	0.04572	8319.38
18	13.46	14939	Aldrin	V	0.00137	4373.26
19	13.77	83571		в	0.08357	22150.76
20	13.89	579595		V	0.57959	150168.44
21	14.05	18995		V	0.01899	5399.64
22	14.12	44462		V	0.04446	12164.17
23	14.20	41793		V	0.04179	12700.27
24	14.29	167832		V	0.16783	36485.04
25	14.48	76044		V	0.07604	23068.17
26	14.61	188265		В	0.18827	41771.53
27	14.84	327383		в	0.32738	99769.34
28	15.00	238956	Hept, epoxide	V	0.00770	36950.20
29	15.09	264130		V	0.26413	64920.81
30	15.27	2188010	gamma chlordane	В	0.06723	584535.03
31	15.55	3649934	alpha chlordane	В	0.11965	798381.01
32	15.68	22538		V	0.02254	8137.44
33	15.94	38156	Endosulfan I	В	0.00145	11318.83
34	16.04	139100		V	0.13910	38478.06
35	16.12	30153		V	0.03015	11052.63
36	16.26	105241		В	0.10524	25656.98
37	16.46	40918	Dieldrin	В	0.00223	11538.93
38	16.58	109371		V	0.10937	24435.06
39	16.74	20412		В	0.02041	6213.36
40	16.81	31556		V	0.03156	8341.21
41	16.89	37161	Endrin	V	0.00269	10910.89
42	17.01	453257	4,4'-DDD	V	0.02021	112359.90
43	17.27	502860		В	0.50286	132663.47
44	17.44	15630	Endosulfan II	V	9.23e-04	4159.47
45	17.81	119515		в	0.11952	29058.53
46	17.90	14986		E	0.01499	4645.78
47	18.21	33908		В	0.03391	6757.89
48	18.38	5791	Endrin aldehyde	V	5.45e-04	1674.59
49	18.51	4526		В	0.00453	1362.69
50	19.11	3457		В	0.00346	1246.43
51	19.27	24193	Endo. Sulfate	В	0.00408	4458.45
52	19.39	36464		V	0.03646	7290.66
53	19.53	6126		V	0.00613	1775.44
54	22.12	9172	Decachlorobiphenyl	В	-4.5e-04	594.06

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4.31982 2.98e+06

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Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98806	Date	: 09/29/2009 07:49:53
Operator	: tchrom	Sample Name	: 9090313
Sample Number	:	Study	: TOX
AutoSampler	: BUILT-IN	Rack/Vial	: 1/31
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Samoling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ut	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 21:09:07	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46231.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46231.rst Inst Method : h:\turbo6\6890-06\6890-612-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46231.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46231.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46231.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	3304		В	0.00330	1280.94
2	11.48	17518		В	0.01752	5509.25
3	12.59	10983		В	0.01098	3331.46
4	13.17	3333		В	0.00333	1123.79
5	13.39	11029		В	0.01103	3440.06
6	13.54	3227	Aldrin	В	0.00102	1025.93
7	13.80	2908		В	0.00291	1095.22
8	13.99	7477		В	0.00748	2477.32
9	14.43	5986		в	0.00599	2062.47
10	14.60	16679		В	0.01668	3667.98
11	14.68	23494		V	0.02349	6011.28
12	14.76	40501		V	0.04050	8783.05
13	14.87	20569		V	0.02057	4998.97
14	15.01	76272	Hept. epoxíde	В	0.00243	16817.58
15	15.27	9508	gamma chlordane	В	0.00116	2696.25
16	15.56	71836	alpha chlordane	В	0.00273	13446.28

09/29/2009 07:49:53 Result: H:\TURBO6\6890-06\6-SEQ46\6A46231.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	15 68	15810		в	0.01581	3929.02
18	15.80	15503	4 4'-DDE	v	0.00174	3432.39
19	15.89	35390	Endosulfan	v	0.00135	6425.60
20	16.00	11660	LindoSullari	v	0.00100	2562 71
20	16 21	16408		Ř	0.01641	4439 04
22	16 31	48699		v	0.01041	13994 46
23	16.36	45717		v	0.04070	14787 07
20	16.00	97017	Dieldrin	v	0.0400	23846.04
24	16.62	187244	Diciuliii	Ň	0.00400	35844 08
20	16.82	360757		v	0.10724	5/881 /9
20	16.02	2/3/56	Endrin	Ň	0.00070	30006.06
21	17 14	60100		Ň	0.01010	10357 76
20	17.14	120467	4,4-000	v	0.00397	28487.00
29	17.21	204479	Endoculfon II	Ň	0.13347	12382.04
21	17.41	294470	Endosulian n	v	0.01207	95152 22
20	17.00	262501		Ň	0.47100	01103.00
JZ 22	17.00	302301	4,4-001		0.01732	40246 24
22	17.00	40091			0.04309	10310.24
34	10.02	30209		÷	0.03027	10270.10
30	10.02	1/0430		v.	0.1/044	77129 20
30	10.11	203724		Ň	0,20072	05054.00
31	10.10	504005	Carlein aldahurda	Ň	0.30401	00004.92
30	10.29	332223	Endrin aldenyde	Ň	0.03237	04401.73
39	10.41	077445			0.17301	30393.40
40	10.07	21/410		V V	0.2/741	00024.33
41	10.04	315047	Math as use blass	Ň	0.31305	00000.29
42	10.78	204449	Methoxychior		0.01885	44190.91
43	10.07	419000		V.	0.41905	03991.03
44	19.07	544410		. <u>.</u>	0.54441	101299.33
45	19.18	//41/8		V.	0.77418	89220.07
40	19.43	162713		V.	0.16271	41102.14
47	19.57	245320		N.	0.24532	45370.51
48	19.68	118352		<u> </u>	0.11835	23143.30
49	19.85	321601	Endrin Ketone	V.	0.01487	55823.47
50	20.05	133051		V.	0.13305	21537.26
51	20.17	159345		<u> </u>	0.15934	37180.08
52	20.30	60158		V.	0.06016	15205.02
53	20.35	50749		V.	0.05075	13802.06
54	20.44	25089		V	0.02509	7095.33
55	20.57	11475		V	0.01148	2817.85
56	20.70	11857		V	0.01186	2717.99
57	20.95	9034		B	0.00903	2457.49
58	21.08	11080		V	0.01108	2996.02
59	21.67	7266		В	0.00727	1946.46
60	22.09	24247	Decachlorobiphenyl	В	2.23e-04	1506.98
		8260917			6.02430	1.59e+06

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:50:57
Reprocess Number	: buf2048: 98820		
Operator	: tchrom	Sample Name	: 9090864-CCV2
Sample Number		Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/38
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29,47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reiect	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 01:20:31	Cycle	: 12

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46238.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46238.rst

Result File : H:\TURBO6\6890-06\6890-06\6-SEQ46\6A46238.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46238.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46238.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46238.rst Report Format File: h:\turbo6\6890-06\06\d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time	ΒL	Area	Component	NG	Height	%D	RT Window -	Relative
[min]		[uV-sec]	Name	CONC.	[µV]	0.05ng		
10.77	BB	1908096	alpha-BHC	0.04947	628936.28	-1.1	10.72 -	10.82
11.61	BB	1754518	gamma-BHC	0.04983	560592.22	-0.3	11.56 -	11.66
11.86	BB	723063	beta-BHC	0.05047	225687.18	0.9	.11.81 -	11.91
12.30	BB	1798075	delta-BHC	0.05038	568451.56	0.8	12.25 -	12.35
12.82	BB	1749144	Heptachlor	0.05249	553568.28	5.0	12.77 -	12.87
13.55	BB	1643953	Aldrin	0.05083	515127.99	1.7	13.50 -	13.60
15.00	BB	1551779	Hept. epoxide	0.05027	466994.43	0.5	14.95 -	15.05
15.28	BB	1584345	gamma chiordane	0.04892	487775.99	-2.2	15.23 -	15.33
15.58	BB	1473215	alpha chlordane	0.04852	455716.79	-3.0	15.53 -	15.63
15.79	ΒV	1494080	4,4'-DDE	0.04883	475475.51	-2.3	15.74 -	15.84
15.89	VB	1387831	Endosulfan I	0.04979	406648.60	-0.4	15.84 -	15.94
16.43	BB	1513075	Dieldrin	0.04868	445395.68	-2.6	16.38 -	16.48
16.94	BB	1267203	Endrin	0.04734	366201.62	-5.3	16.89 -	16.99
17.11	BB	1157462	4,4'-DDD	0.04997	360663.18	-0.1	17.06 -	17.16
17.44	BB	1207370	Endosulfan II	0.04858	338542.85	-2.8	17.39 -	17.49
17.71	BB	1085994	4,4'-DDT	0.04668	327347.03	-6.6	17.66 -	17.76

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09/29/2009 07:50:57 Result: H:\TURBO6\6890-06\6-SEQ46\6A46238.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - F	Relative
18.36	BB	799678	Endrin aldehvde	0.04854	218845.36	-2.9	18.31 -	18.41
18.76	BB	538769	Methoxychlor	0.04776	162525.05	-4.5	18.71 -	18.81
19.30	BB	974771	Endo, Sulfate	0.05269	268479.80	5.4	19.25 -	19.35
19.90	BB	1212544	Endrin ketone	0.05046	319196.77	0.9	19.85 -	19.95
		<u> </u>						
		26824966		0.99049	8.15e+06			
Missing Component Report Component Expected Retention (Calibration File)								

Tetrachloro-m-xylene9.178Decachlorobiphenyl22.082

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Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98822	Date	: 09/29/2009 07:51:04
Operator	: tchrom	Sample Name	: 9090254-CCV2
Sample Number	•	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/39
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 01:56:31	Cycle	: 13

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46239.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46239.rst

Result File : H:\1URB06\6890-06\6-SEQ46\6A46239.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURB06\6890-06\6-SEQ46\6A46239.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURB06\6890-06\6-SEQ46\6A46239.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURB06\6890-06\6-SEQ46\6A46239.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURB06\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window	•	Relative
9.19	BB	1211949	Tetrachloro-m-xylene	0.04763	393599.68	-4.7	9.14	-	9.24
15.27	BB	4525	gamma chlordane	0.00101	1536.43	-98:0	15.22	-	15.32
15.92	BB	20225	Endosulfan I	8.06e-04	362.67	-98.4	15.87	-	15.97
17.73	BB	11636	4,4'-DDT	0.00307	498.27	-93.9	17.68	-	17.78
22.10	BB	1108317	Decachlorobiphenyl	0.04859	267131.08	-2.8	22.05	-	22.15
		2356653		0.10112	663128.13				

9/29/09

Missing Component Report

Component Expected Retention (Calibration File)

alpha-BHC	10.743
gamma-BHC	11.590
beta-BHC	11.839
delta-BHC	12.280
Heptachlor	12.799
Aldrin	13.529
Hept. epoxide	14.977
alpha chlordane	15.566
4,4'-DDE	15.768

9/29/00

Page 1 of 2

Software Version	: 6.2.1.0.104:0104	Date : 09/29/2009 07:51:47	,
Reprocess Number	: buf2048: 98836		
Operator	: tchrom	Sample Name : 9090864-CCV3	
Sample Number	:	Study : 8081	
AutoSampler	: BUILT-IN	Rack/Vial : 1/46	
Instrument Name	: HP6890-06	Channel : A	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.45 min	
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject : 3000.000000	
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	: 09/29/2009 06:08:05	Cycle : 20	

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46246.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46246.rst

Inst Method : h:\turbo6\6890-06\6a90-0

Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time	BL	Area	Component	NG	Height	%D	RT Window -	Relative
[[[[[[]]]]		[uv-sec]			[v u]	0.05iig	· · · · · · · · · · · ·	
10.75	BΒ	1952376	alpha-BHC	0.05059	641676.50	1.2	10.70 -	10.80
11.60	BB	1806332	gamma-BHC	0.05126	577977.43	2.5	11.55 -	11.65
11.85	BB	747070	beta-BHC	0.05217	232793.60	4.3	11.80 -	11.90
12.29	BB	1869672	delta-BHC	0.05231	595750.26	4.6	12.24 -	12.34
12.81	BΒ	1804754	Heptachlor	0.05415	568045.71	8.3	12.76 -	12.86
13.54	BB	1726554	Aldrin	0.05334	543540.43	6.7	13.49 -	13.59
14.98	BB	1630838	Hept. epoxide	0.05283	488449.02	5.7	14.93 -	15.03
15.27	BB	1671718	gamma chlordane	0.05157	512890.84	3.1	15.22 -	15.32
15.57	BΒ	1579055	alpha chlordane	0.05198	481268.75	4.0	15.52 -	15.62
15.77	ΒV	1567181	4,4'-DDE	0.05116	500963.58	2.3	15.72 -	15.82
15.88	VB	1451713	Endosulfan I	0.05208	428842.01	4.2	15.83 -	15.93
16.41	BB	1595130	Dieldrin	0.05127	473800.34	2.5	16.36 -	16.46
16.93	BΒ	1341812	Endrin	0.05005	383573.79	0.1	16.88 -	16.98
17.09	BB	1275204	4,4'-DDD	0.05495	397268.77	9.9	17.04 -	17.14
17.42	BΒ	1270135	Endosulfan II	0.05109	355741.62	2.2	17.37 -	17.47
17.70	VB	1001200	4,4'-DDT	0.04324	305491.21	-13.5	17.65 -	17.75
18.34	BΒ	836184	Endrin aldehyde	0.05074	226388.64	1.5	18.29 -	18.39

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09/29/2009 07:51:47 Result: H:\TURBO6\6890-06\6-SEQ46\6A46246.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - f	Relative
18.75	BB	524737	Methoxychlor	0.04654	158010.02	-6.9	18.70 -	18.80
19.29	BB	1036178	Endo. Sulfate	0.05583	285398.69	11.7	19.24 -	19.34
19.89	BΒ	1267676	Endrin ketone	0.05266	331370.16	5.3	19.84 -	19.94
22.07	BB	19301	Decachlorobiphenyl	2.44e-06	1278.74	-100.0	22.02 -	22.12
		27974821		1.02981	8.49e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene

9.178

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Software Version	: 6.2.1.0.104:0104	Date : 09/29/2009 07:51:52	
Reprocess Number	: buf2048: 98838		
Operator	: tchrom	Sample Name : 9090254-CCV3	
Sample Number	:	Study : SURR	
AutoSampler	: BUILT-IN	Rack/Vial : 1/47	
Instrument Name	: HP6890-06	Channel : A	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.49 min	
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject : 3000.000000	
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	: 09/29/2009 06:44:03	Cycle : 21	

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46247.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46247.rst

Result File : H:\1URBO6\6890-06\6890-06\6-SEQ46\6A46247.rst Inst Method : h:\turbo6\6890-06\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46247.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46247.rst Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46247.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window	- 1	Relative	•
9.18 15.26 15.78 22.08	BB BB BB BB	1224055 3184 10194 1166423	Tetrachloro-m-xylene gamma chlordane 4,4'-DDE Decachlorobiphenyl	0.04810 9.69e-04 0.00157 0.05119	401131.45 1181.99 521.64 281967.58	-3.8 -98_1 -96_9 2.4	9.13 15.21 15.73 22.03	- - -	9.23 15.31 15.83 22.13	9
		2403855		0.10183	684802.65					

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Page 1 of 2

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98840	Date	: 09/29/2009 07:51:56
Operator	: tchrom	Sample Name	: 9091475-CCV3
Sample Number	:	Study	: PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/48
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.48 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 07:20:06	Cycle	: 22

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46248.raw <Modified>



Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46248.raw

Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46248.rst

Calib Method : h:\turbo6\6890-06\6a-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6A46248.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.76	3602		В	0.00360	1408.76
2	9.18	467796	Tetrachloro-m-xylene	В	0.01895	153134.66
3	10.75	305252	alpha-BHC	В	0.00879	100319.92
4	11.60	306313	gamma-BHC	в	0.00966	97304.47
5	11.85	130095	beta-BHC	В	0.00860	39881.08
6	15.26	6056	gamma chlordane	В	0.00106	2014.84
7	15.78	21098	4,4'-DDE	В	0.00192	3338.58
8	16.72	24904		В	0.02490	6881.19
9	16.93	1253537	Endrin	В	0.04685	359397.28
10	17.10	111668	4,4'-DDD	V	0.00577	33832.90
11	17.70	2202824	4,4'-DDT	В	0.09201	669073.90
12	18.03	4130		В	0.00413	1320.79
13	18.34	26575	Endrin aldehyde	В	0.00180	7550.26
14	18.60	12164	-	В	0.01216	3076.48



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09/29/2009 07:51:56 Result: H:\TURBO6\6890-06\6-SEQ46\6A46248.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL.	NG CONCENTRATION	Height [µV]
15	18.75	2751829	Methoxychlor	v	0.23909	839839.50
16	18.93	9973	,	В	0.00997	3159.06
17	19.89	54595	Endrin ketone	В	0.00421	14448.85
18	22.08	467335	Decachlorobiphenyl	В	0.01999	112706.77
		8159746			0.51347	2.45e+06

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:49:18
Reprocess Number	: buf2048: 98799		
Operator	: tchrom	Sample Name	: 9091475-CCV1
Sample Number	:	Study	: PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/27
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.00000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 18:44:59	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46227.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46227.rst



Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46227.raw

Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46227.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46227.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.09	1819	<u></u>	В	0.00182	761.70
2	8.28	2134		В	0.00213	803.01
3	8.61	1516		В	0.00152	597.24
4	9.10	1685		В	0.00168	733.12
7	10.53	195127	Tetrachloro-m-xylene (В	0.02103	60272.68
8	10.93	1875		В	0.00187	594.10
12	12.30	129792	alpha-BHC (2C)	V	0.01053	40650.13
15	13.25	133835	gamma-BHC (2C)	В	0.01079	40121.15
16	13.47	59391	beta-BHC (2C)	В	0.00937	17478.17
17	15.34	3267		В	0.00327	888.67
19	17.68	7037	4,4'-DDE (2C)	В	0.00187	2106.67
20	18.77	482583	Endrin (2C)	В	0.05275	128479.76
21	18.96	9746	4,4'-DDD (2C)	V	0.00228	2810.45
22	19.51	2972	, , ,	В	0.00297	830.08
23	19.65	855909	4,4'-DDT (2C)	V	0.11006	248443.78
24	19.95	7018	Endrin aldehyde (2C)	В	1.72e-04	1843.43



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09/29/2009 07:49:18 Result: H:\TURBO6\6890-06\6-SEQ46\6b46227.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
26	20.62	2394	Endo. Sulfate (2C)	В	0.00227	726.57
27	20.76	2621		в	0.00262	626.35
28	21.09	1036770	Methoxychlor (2C)	В	0.27986	290772.80
29	21.37	4450		в	0.00445	1276.39
30	21.78	13197	Endrin ketone (2C)	В	0.00264	3071.03
31	23.60	11662		В	0.01166	272.07
32	25.22	204836	Decachlorobiphenyl (2C	В	0.02207	34783.70
		3171634			0.55969	878943.07

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:49:28
Reprocess Number	: buf2048: 98801		
Operator	: tchrom	Sample Name	: 9090864-CCV1
Sample Number		Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/28
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 19:21:06	Cycle	: 2

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46228.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46228.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46228.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46228.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)sur(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46228.rst Report Format File: h:\turbo6\6890-06\06%d.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6b-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.29	BB	742291	alpha-BHC (2C)	0.04996	235235.86	-0.1	12.24 -	12.34
13.25	BΒ	701114	gamma-BHC (2C)	0.05056	214640.14	1.1	13.20 -	13.30
13.47	BB	302948	beta-BHC (2C)	0.05201	89862.34	4.0	13.42 -	13.52
14.22	VB	698480	delta-BHC (2C)	0.05054	211600.91	1.1	14.17 -	14.27
14.40	BB	669889	Heptachlor (2C)	0.05104	202540.59	2.1	14.35 -	14.45
15.20	ΒV	621550	Aldrin (2C)	0.05117	186139.17	2.3	15.15 -	15.25
16.55	BB	591286	Hept. epoxide (2C)	0.05207	167039.65	4.1	16.50 -	16.60
17.00	BB	584926	gamma chlordane (2C)	0.05179	169043.49	3.6	16.95 -	17.05
17.33	ΒV	554688	alpha chlordane (2C)	0.05193	157234.68	3.9	17.28 -	17.38
17.47	VB	527097	Endosulfan I (2C)	0.05201	146829.66	4.0	17.42 -	17.52
17.68	BΒ	542000	4,4'-DDE (2C)	0.05115	165485.32	2.3	17.63 -	17.73
18.08	BB	563431	Dieldrin (2C)	0.05122	157004.64	2.4	18.03 -	18.13
18.77	BB	460943	Endrin (2C)	0.05045	123129.88	0.9	18.72 -	18.82
18.95	BΒ	425509	4,4'-DDD (2C)	0.05138	124103.13	2.8	18.90 -	19.00
19.22	BB	463052	Endosulfan II (2C)	0.05180	122287.62	3.6	19.17 -	19.27
19.65	BB	384021	4,4'-DDT (2C)	0.05069	111209.71	1.4	19.60 -	19.70

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09/29/2009 07:49:28 Result: H:\TURBO6\6890-06\6-SEQ46\6b46228.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
19.95	ΒV	317731	Endrin aldehyde (2C)	0.05215	81840.46	4.3	19.90 -	20.00
20.56	BB	359313	Endo. Sulfate (2C)	0.05183	92967.22	3.7	20.51 -	20.61
21.09	BB	197656	Methoxychlor (2C)	0.05343	54351.20	6.9	21.04 -	21.14
21.78	BB	455746	Endrin ketone (2C)	0.05243	102703.49	4.9	21.73 -	21.83
		10163672		1.02959	2.92e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene (2C)	10.530
Decachlorobiphenyl (2C)	25.216

129/09 min

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Software Version	: 6.2.1.0.104:0104	Date : 09/29/2009 09:59:42	
Reprocess Number	: buf2048: 98865		
Operator	: tchrom	Sample Name : 9090254-CCV1	
Sample Number	:	Study : SURR	
AutoSampler	: BUILT-IN	Rack/Vial : 1/29	
Instrument Name	: HP6890-06	Channel : B	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.46 min	
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject : 3000.000000	
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	: 09/28/2009 19:57:04	Cycle : 1	

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46229.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46229.rst

Result File : H:\1URBO6\6890-06\65EQ46\6b46229.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46229.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46229.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46229.rst Report Format File: h:\turbo6\6890-06\06\deltag0-06\06\deltag1.pt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window	-	Relative
10.54 25.22	BB BB	456695 450019	Tetrachloro-m-xylene (Decachlorobiphenyl (2C	0.04838 0.05103	142374.47 75954.45	-3.2 2.1	10.49 25.17	-	10.59 25.27
		906714		0.09941	218328.92				

Missina	Component	Report
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Expected Retention (Calibration File) Component

alpha-BHC (2C)	12.289
gamma-BHC (2C)	13.247
beta-BHC (2C)	13.465
delta-BHC (2C)	14.214
Heptachlor (2C)	14.398
Aldrin (2C)	15.200
Hept. epoxide (2C)	16.551
gamma chlordane (2C)	16.989
alpha chlordane (2C)	17.322
Endosulfan I (2C)	17.458
4,4'-DDE (2C)	17.672
Dieldrin (2C)	18.075

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:49:48
Reprocess Number Operator Sample Number AutoSampler Instrument Name Instrument Serial #	: but2048: 98805 : tchrom : : BUILT-IN : HP6890-06 : CN10520010	Sample Name Study Rack/Vial Channel A/D mV Range	: 9090308 : TECH : 1/30 : B : 1000 : 20.47 min
Delay Time Sampling Rate Sample Volume Sample Amount Data Acquisition Time	: 0.00 min : 5.0000 pts/s : 1.00000 ul : 1.0000 : 09/28/2009 20:32:59	Area Reject Dilution Factor Cycle	: 1500.000000 : 1.00 : 4

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46230.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46230.rst Inst Method : h:\turbo6\6890-06\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46230.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46230.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)sur(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46230.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	1996		В	0.00200	755.41
2	8.35	2146		V	0.00215	750.16
3	10.64	60874	Tetrachloro-m-xylene (В	0.00699	18840.01
4	12.17	10927		В	0.01093	3316.47
5	12.40	8473		В	0.00847	2533.09
6	13.02	23676		В	0.02368	7017.47
7	13.13	5632		V	0.00563	1514.90
8	13.31	4757	gamma-BHC (2C)	В	0.00174	1319.18
9	13.58	16355		В	0.01635	5095.77
10	13.73	18190		В	0.01819	5006.89
11	13.82	7809		V	0.00781	2134.59
12	13.95	27322		V	0.02732	8142.32
13	14.03	181877		V	0.18188	53158.08
14	14.40	374176	Heptachlor (2C)	В	0.02879	109089.84
15	14.67	24934		В	0.02493	7387.63
16	14.77	32202		V	0.03220	7312.60

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09/29/2009 07:49:48 Result: H:\TURBO6\6890-06\6-SEQ46\6b46230.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	14.95	4778		в	0.00478	1347.08
18	15.09	6315		В	0.00631	2074.34
19	15.34	52014		В	0.05201	14521.06
20	15.56	211000		В	0.21100	54746.73
21	15.66	7487		Ε	0.00749	3719.54
22	15.70	18132		V	0.01813	5084.14
23	15.86	35164		V	0.03516	7191.01
24	15.96	28055		V	0.02806	7266.14
25	16.04	83115		V	0.08311	20467.38
26	16.18	34863		V	0.03486	9662.61
27	16.31	5754		В	0.00575	1666.35
28	16.47	159370		V	0.15937	39159.08
29	16.62	12581	Hept. epoxide (2C)	V	0.00114	3044.93
30	16.74	144372		V	0.14437	32572.16
31	16.84	44604		V	0.04460	8838.21
32	17.00	754850	gamma chlordane (2C)	v	0.06671	217013.54
33	17.22	806497		V	0.80650	205657.53
34	17.33	533776	alpha chlordane (2C)	V	0.04998	151421.55
35	17.53	4885	Endosulfan I (2C)	В	4.78e-04	1192.52
36	17.65	28774	4,4'-DDE (2C)	V	0.00387	5441.26
37	17.88	52603		В	0.05260	13609.79
38	18.11	54689	Dieldrin (2C)	V	0.00579	9220.65
39	18.31	9565		В	0.00956	3020.50
40	18.38	18249		V	0.01825	4585.98
41	18.49	27393		V	0.02739	4835.04
42	18.59	2877		V	0.00288	889.62
43	18.70	29659	Endrin (2C)	V	0.00445	5295.96
44	18.84	15787		V	0.01579	3923.04
45	18.94	165149	4,4'-DDD (2C)	V.	0.02063	43692.29
46	19.09	28493		V	0.02849	6268.33
47	19.34	179982		v N	0.17998	43242.34
48	19.53	8352		В	0.00835	2140.42
49	19.61	6300	4,4°-DDT (2C)	V V	0.00317	1932.45
50	20.05	35746		В	0.03575	9212.83
51	20.21	2486		B	0.00249	690.35
52	20.32	9555		v N	0.00955	2430.08
53	20.63	1667	Endo. Sulfate (2C)	В	0.00216	520.51
54	20.78	3835	14.00	В	0.00383	1033.75
55	21.03	4292	Methoxychior (2C)	В	0.00123	1203.30
50	21.07	12828	Endrin katana (DC)	B	0.01203	1000.04
5/	21.02	0121	Enunn Kelone (20)		0.00184	620 49
ວຽ	22.10	2395		D	0.00240	020.40
		4455751			2.61219	1.19e+06

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:49:58
Reprocess Number	: buf2048: 98807		
Operator	: tchrom	Sample Name	: 9090313
Sample Number	•	Study	: TOX
AutoSampler	: BUILT-IN	Rack/Vial	: 1/31
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.49 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/28/2009 21:09:07	Cycle	: 5

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46231.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46231.rst

Result File : H:\1URBO6\6890-06\6890-06\6-SEQ46\6b46231.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46231.rsw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46231.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46231.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.29	2064	<u>,</u>	В	0.00206	754.83
2	13.08	8111		В	0.00811	2403.01
3	14.07	4254		В	0.00425	1281.16
4	14.22	1872	delta-BHC (2C)	В	0.00219	534.43
5	14.84	4376		В	0.00438	1298.87
6	15.34	3772		В	0.00377	1030.46
7	15.44	2726		V	0.00273	926.09
9	16.21	16884		В	0.01688	4519.16
10	16.40	6719		V	0.00672	1795.56
11	16.53	6084	Hept. epoxide (2C)	В	5.67e-04	1699.36
12	16.62	24673	,	V	0.02467	6342.86
15	17.15	10157		В	0.01016	2819.05
16	17.44	2429	Endosulfan I (2C)	В	2.36e-04	829.73
18	17.70	10814	4,4'-DDE (2C)	В	0.00222	1751.69
19	17.92	8131		В	0.00813	2413.35
20	18.03	65537	Dieldrin (2C)	V	0.00675	13632.34

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09/29/2009 07:49:58 Result: H:\TURBO6\6890-06\6-SEQ46\6b46231.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL		Height [µV]
21	18.16	7686		v	0.00769	1620.65
22	18.29	47934		v	0.04793	11498.61
23	18.43	48206		Ý	0.04821	10209.88
24	18.54	32485		v	0.03248	9080.82
25	18.63	141184		Ý	0.14118	23046.30
26	18.78	38924	Endrin (2C)	v	0.00544	7362.65
27	18.86	19880		V	0.01988	5378.66
28	18.97	57785	4.4'-DDD (2C)	v	0.00796	9397.31
29	19.12	60315	, , , , , , , , , , , , , , , , , , , ,	V	0.06032	12962.92
30	19.20	59632	Endosulfan II (2C)	v	0.00666	12519.00
31	19.31	41047		v	0.04105	10963.53
32	19.41	231926		v	0.23193	49225.50
33	19.58	111514	4,4'-DDT (2C)	V	0.01640	18690.78
34	19.83	101861	·,· - ()	v	0.10186	17783.06
35	19.98	251895	Endrin aldehyde (2C)	V	0.04114	45682.23
36	20.17	180040	, , ,	v	0.18004	26783.16
37	20.29	62153		V	0.06215	10180.77
38	20.47	109720		V	0.10972	17544.97
39	20.56	115564	Endo, Sulfate (2C)	V	0.01798	24684.26
40	20.77	238652		V	0.23865	19734.95
41	21.00	82322		V	0.08232	14153.78
42	21.11	191746	Methoxychlor (2C)	V	0.05183	34996.37
43	21.23	217880		V	0.21788	35571.50
44	21.37	69742		V	0.06974	14873.90
45	21.57	67864		V	0.06786	12726.82
46	21.68	99860		V	0.09986	13589.31
47	21.92	77307		V	0.07731	11967.89
48	22.05	87968		V	0.08797	15606.69
49	22.23	31006		V	0.03101	4368.24
50	22.35	29326		V	0.02933	5962.25
51	22.48	59145		V	0.05915	9291.17
52	22.63	19924		V	0.01992	3906.07
53	22.78	24834		V	0.02483	3083.06
54	22.96	3494		V	0.00349	869.42
55	23.49	8095		В	0.00809	946.77
		3207519			2.45310	570295.17

9/29/09 Mm

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Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:51:00
Reprocess Number	: but2048: 98821	Cample Name	+ 0000864 CCV2
Operator	: tonrom	Sample Name	. 9090604-CCVZ
Sample Number		Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/38
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.47 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 01:20:31	Cycle	: 12

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46238.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46238.rst

 Result File:
 In:/turbo6\6890-06\6890-612-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46238.raw

 Proc Method :
 h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46238.rst

 Calib Method :
 h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46238.rst

 Report Format File:
 h:\turbo6\6890-06\06\06\06\075_C

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL.	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.31	BB	737235	alpha-BHC (2C)	0.04963	234233.92	-0.7	12.26 -	12.36
13.26	BB	696930	gamma-BHC (2C)	0.05026	212794.04	0.5	13.21 -	13.31
13.48	BB	299918	beta-BHC (2C)	0.05148	89083.85	3.0	13.43 -	13.53
14.23	BΒ	695765	delta-BHC (2C)	0.05035	210459.39	0.7	14.18 -	14.28
14.41	BB	671406	Heptachlor (2C)	0.05115	203688.55	2.3	14.36 -	14.46
15.22	ΒV	603906	Aldrin (2C)	0.04975	180482.63	-0.5	15.17 -	15.27
16.57	BΒ	568588	Hept. epoxide (2C)	0.05007	161734.20	0.1	16.52 -	16.62
17.01	BB	554898	gamma chlordane (2C)	0.04916	160454.92	-1.7	16.96 -	17.06
17.34	ΒV	527731	alpha chlordane (2C)	0.04941	150901.37	-1.2	17.29 -	17.39
17.48	VB	503355	Endosulfan I (2C)	0.04967	142830.04	-0.7	17.43 -	17.53
17.69	BΒ	516069	4,4'-DDE (2C)	0.04876	158044.50	-2.5	17.64 -	17.74
18.10	BΒ	535597	Dieldrin (2C)	0.04874	150114.81	-2.5	18.05 -	18.15
18.78	BΒ	422207	Endrin (2C)	0.04631	114123.83	-7.4	18.73 -	18.83
18.97	BΒ	410519	4,4'-DDD (2C)	0.04961	120624.40	-0.8	18.92 -	19.02
19.23	ΒB	440553	Endosulfan II (2C)	0.04928	118293.06	-1.4	19.18 -	19.28
19.66	BB	363361	4,4'-DDT (2C)	0.04809	105660.82	-3.8	19.61 -	19.71

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09/29/2009 07:51:00 Result: H:\TURBO6\6890-06\6-SEQ46\6b46238.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - F	Relative
19.96	ΒV	298811	Endrin aldehyde (2C)	0.04898	77905.25	-2.0	19.91 -	20.01
20.57	BB	375832	Endo. Sulfate (2C)	0.05413	98352.91	8.3	20.52 -	20.62
21.10	BΒ	185126	Methoxychlor (2C)	0.05004	51480.42	0.1	21.05 -	21.15
21.80	BΒ	454524	Endrin ketone (2C)	0.05229	103451.07	4.6	21.75 -	21.85
		9862332		0.99716	2.84e+06			

Missing Component Report Component Expected Retention (Calibration File)

Tetrachloro-m-xylene (2C)	10.530
Decachlorobiphenyl (2C)	25.216

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Page 1 of 2

Software Version	: 6.2.1.0.104:0104	Date	: 09/29/2009 07:51:08
Reprocess Number	: buf2048: 98823		
Operator	: tchrom	Sample Name	: 9090254-CCV2
Sample Number	:	Study	: SURR
AutoSampler	: BUILT-IN	Rack/Vial	: 1/39
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delav Time	: 0.00 min	End Time	: 29.50 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reiect	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 01:56:31	Cycle	: 13

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46239.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46239.rst

Result File : H:\1URB06\6890-06\6890-6 12-5-08-ins from H:\TURB06\6890-06\6-SEQ46\6b46239.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURB06\6890-06\6-SEQ46\6b46239.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURB06\6890-06\6-SEQ46\6b46239.rst Report Format File: h:\turbo6\6890-06\06\deltage.pst Sequence File : H:\TURB06\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - Re	lative
10.54 18.75 25.24	BB BB BB	468343 4333 427410	Tetrachloro-m-xylene (Endrin (2C) Decachlorobiphenyl (2C	0.04960 0.00175 0.04836	146657.12 100.67 72990.92	-0.8 - <u>96 5</u> -3.3	10.49 - 18.70 - 25.19 -	10.59 18.80 25.29
		900086		0.09971	219748.71			

Missing Component Report

Component Expected Retention (Calibration File)

alpha-BHC (2C)	12.289
gamma-BHC (2C)	13.247
beta-BHC (2C)	13.465
delta-BHC (2C)	14.214
Heptachlor (2C)	14.398
Aldrin (2C)	15.200
Hept. epoxide (2C)	16.551
gamma chlordane (2C)	16.989
alpha chlordane (2C)	17.322
Endosulfan I (2C)	17.458
4,4'-DDE (2C)	17.672

Page 1 of 2

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98837	Date	: 09/29/2009 07:51:50
Operator	: tchrom	Sample Name	: 9090864-CCV3
Sample Number	:	Study	: 8081
AutoSampler	: BUILT-IN	Rack/Vial	: 1/46
Instrument Name	: HP6890-06	Channel	: B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 3000.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 06:08:05	Cycle	: 20

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46246.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46246.rst

 Result File
 In: \UCDO0/0690-06/6890-06/6890-6

 Inst Method
 : h:\turbo6\6890-06\6890-6
 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46246.raw

 Proc Method
 : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46246.rst

 Calib Method
 : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46246.rst

 Report Format File
 h:\turbo6\6890-06\06\%d.rpt

 Sequence File
 H:\TURBO6\6890-06\6-SEQ46\6b-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window -	Relative
12.29	BB	781566	alpha-BHC (2C)	0.05248	249168.93	5.0	12.24 -	12.34
13.25	BB	732944	gamma-BHC (2C)	0.05279	223512.00	5.6	13.20 -	13.30
13.47	BB	314802	beta-BHC (2C)	0.05408	93487.27	8.2	13.42 -	13.52
14.22	BB	733420	delta-BHC (2C)	0.05296	222071.05	5.9	14.17 -	14.27
14.40	BB	697633	Heptachlor (2C)	0.05313	210386.09	6.3	14.35 -	14.45
15.20	BV	637045	Aldrin (2C)	0.05242	191049.85	4.8	15.15 -	15.25
16.56	BB	597868	Hept. epoxide (2C)	0.05265	170863.94	5.3	16.51 -	16.61
17.00	BB	586093	gamma chlordane (2C)	0.05190	169660.88	3.8	16.95 -	17.05
17.33	ΒV	559409	alpha chlordane (2C)	0.05237	160151.27	4.7	17.28 -	17.38
17.47	VB	531802	Endosulfan I (2C)	0.05247	148606.30	4.9	17.42 -	17.52
17.68	BB	546256	4,4'-DDE (2C)	0.05154	167287.65	3.1	17.63 -	17.73
18.08	BΒ	568052	Dieldrin (2C)	0.05163	157906.22	3.3	18.03 -	18.13
18.77	BB	460269	Endrin (2C)	0.05037	124657.86	0.7	18.72 -	18.82
18.95	BB	458865	4,4'-DDD (2C)	0.05532	134777.16	10.6	18.90 -	19.00
19.22	ΒB	466239	Endosulfan II (2C)	0.05215	125029.43	4.3	19.17 -	19.27
19.65	BB	339939	4,4'-DDT (2C)	0.04514	99014.82	-9.7	19.60 -	19.70

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09/29/2009 07:51:50 Result: H:\TURBO6\6890-06\6-SEQ46\6b46246.rst

Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - F	Relative
19.95	ΒV	316049	Endrin aldehyde (2C)	0.05187	81783.88	3.7	19.90 -	20.00
20.56	BB	401923	Endo, Sulfate (2C)	0.05775	105241.80	15.5	20.51 -	20.61
21.09	BB	183702	Methoxychlor (2C)	0.04966	51195.07	-0.7	21.04 -	21.14
21.78	VB	479754	Endrin ketone (2C)	0.05513	108592.72	10.3	21.73 -	21.83
		10393629		1.04783	2.99e+06			
Missing C	ompo	nent Repo	rt					
Co	mpor	nent .	Expected Retention (C	alibration	File)			

Tetrachloro-m-xylene (2C)	10.530
Decachlorobiphenyl (2C)	25.216

9/29/09

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Software Version	: 6.2.1.0.104:0104	Date : 09/29/2009 07:51:54	
Reprocess Number	: buf2048: 98839		
Operator	: tchrom	Sample Name : 9090254-CCV3	
Sample Number	:	Study : SURR	
AutoSampler	: BUILT-IN	Rack/Vial : 1/47	
Instrument Name	: HP6890-06	Channel : B	
Instrument Serial #	: CN10520010	A/D mV Range : 1000	
Delay Time	: 0.00 min	End Time : 29.49 min	
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.00000 ul	Area Reject : 3000.000000	
Sample Amount	: 1.0000	Dilution Factor : 1.00	
Data Acquisition Time	e : 09/29/2009 06:44:03	Cycle : 21	

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46247.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6b46247.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46247.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46247.rst Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46247.rst Report Format File: h:\turbo6\6890-06\06\06\07.15

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Ret Time [min]	BL	Area [uV-sec]	Component Name	NG CONC.	Height [µV]	%D 0.05ng	RT Window - Rel	ative
10.53 18.75 25.21	BB BB BB	488480 4219 452759	Tetrachloro-m-xylene (Endrin (2C) Decachlorobiphenyl (2C	0.05170 0.00173 0.05135	153943.64 112.87 77310.41	3.4 - 96.5 2.7	10.48 - 1 18.70 - 1 25.16 - 2	0.58 8.80 5.26
		945458		0.10479	231366.93			

Missing Component Report

Expected Retention (Calibration File) Component

······································
12.289
13.247
13.465
14.214
14.398
15.200
16.989
17.322
17.458
17.672
18.075

Page 1 of 2

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98841	Date	: 09/29/2009 07:51:58
Operator Sample Number	: tchrom	Sample Name Study	: 9091475-CCV3 : PEM
AutoSampler	: BUILT-IN	Rack/Vial	: 1/48
Instrument Name	: HP6890-06	Channel	: В
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time Sampling Rate	: 0.00 min : 5.0000 pts/s	End Time	: 29.48 min
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/29/2009 07:20:06	Cycle	: 22

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46248.raw <Modified>



Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46248.raw

Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46248.rst

Calib Method : h:\turbo6\6890-06\6b-pest(09-28-09)surr(9-28-09).mth from H:\TURBO6\6890-06\6-SEQ46\6b46248.rst

Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.28	2302		В	0.00230	848.54
2	10.53	190367	Tetrachloro-m-xylene (В	0.02053	59225.50
3	12.29	125396	alpha-BHC (2C)	В	0.01025	39435.36
4	13.25	128288	gamma-BHC (2C)	В	0.01040	38824.04
5	13.47	57645	beta-BHC (2C)	В	0.00906	16951.75
6	15.34	3770		В	0.00377	1176.35
7	16.66	3318		В	0.00332	110.65
8	17.68	3706	4,4'-DDE (2C)	В	0.00156	1134.79
9	18.76	434272	Endrin (2C)	В	0.04760	115819.83
10	18.95	44107	4,4'-DDD (2C)	V	0.00634	12623.03
11	19.65	737319	4,4'-DDT (2C)	В	0.09514	214683.91
12	19.95	13650	Endrin aldehyde (2C)	В	0.00128	3320.83
14	20.23	1525		В	0.00153	489.73
15	20.62	3555	Endo. Sulfate (2C)	в	0.00243	1027.23

9/29/09 Num 7.7% Endrin 7.7%

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09/29/2009 07:51:58 Result: H:\TURBO6\6890-06\6-SEQ46\6b46248.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [uV]
16 17 18 19	21.09 21.37 21.78 25.21	936220 3827 22425 189160	Methoxychlor (2C) Endrin ketone (2C) Decachlorobiphenyl (2C	B B B B	0.25272 0.00383 0.00368 0.02022	263538.62 1118.92 5152.05 32380.97
		2900853			0.49596	807862.10

Form 1 ORGANIC ANALYSIS DATA SHEET

Blank

8081A

Laboratory: <u>TestAmerica Buffalo</u>				SDG:				
Client: Olin Chlor Alkali Products - Cleveland			veland, TN	d, TN Project:		Olin - Charles Gibson site - NY3A9025AE03759		
Matrix:	<u>Solid</u>	Lab	oratory ID:	<u>9120007</u>	<u>-BLK1</u>	File ID:	<u>6A46193</u>	
Sampled:		Prej	pared:	<u>09/21/09</u>	0.08:00	Analyzed:	<u>09/25/09 12:09</u>	
Solids:		Prej	paration:	<u>3550B C</u>	<u>iC</u>	Initial/Final:	<u>30.2 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>RI92924</u>		Calibration:	<u>R9H1803</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CON	C. (ug/kg)	Q
319-84-6	alpha-BHC				1		1.7	U
319-85-7	beta-BHC				i		U	
319-86-8	delta-BHC				1		1.7	U
58-89-9 gamma-BHC (Lindane)				1	1.7		U	
SYSTEM MONI	TORING COMP	OUND	ADDEI) (ug/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q
Decachlorobiphenyl			6.	62	6.08	92	42 - 146	
Tetrachloro-m-xy	lene		6.	62	5.21	79	37 - 135	<u> </u>

* Values outside of QC limits

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98598	Date	: 09/25/2009 13:14:05
Operator	: tchrom	Sample Name	: 9120007-BLK1
Sample Number	:	Study	:
AutoSampler	: BUILT-IN	Rack/Vial	: 1/93
Instrument Name	: HP6890-06	Channel	: A
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 29.45 min
Sampling Rate	: 5.0000 pts/s		
Sample Volume	: 1.000000 ul	Area Reject	: 1500.000000
Sample Amount	: 1.0000	Dilution Factor	: 1.00
Data Acquisition Time	: 09/25/2009 12:09:51	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46193.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6A46193.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46193.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46193.rst Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46193.rst Report Format File: h:\turbo6\6890-06\6samp.rpt

Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.25	3257		В	0.00326	884.94
2	7.56	6339		В	0.00634	1138.26
3	7.76	3346		В	0.00335	1169.68
4	7.92	100537		В	0.10054	34598.87
5	8.11	11752		В	0.01175	2165.92
6	8.97	34504		В	0.03450	12606.09
7	9.07	2438		V	0.00244	970.02
8	9.36	241249	Tetrachloro-m-xylene	В	0.01574	78471.10
9	9.99	8950	-	В	0.00895	2349.24
10	10.15	5659		В	0.00566	787.43
11	10.62	3322		В	0.00332	1209.12
12	11.00	5152	alpha-BHC	В	0.00363	1191.54
13	11.36	3245	-	В	0.00325	989.69
14	11.46	2613		В	0.00261	1074.40
15	12.13	15189		В	0.01519	4777.51
17	13.50	10748		В	0.01075	3095.67

09/25/2009 13:14:05 Result: H:\TURBO6\6890-06\6-SEQ46\6A46193.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
18	13.85	1785		В	0.00178	666.84
19	14.48	5251		В	0.00525	1955.48
20	14.70	7074		В	0.00707	2151.13
21	14.82	7136		в	0.00714	1990.27
22	14.93	15336		V	0.01534	4112.56
23	15.45	32684	gamma chlordane	В	0.00303	9489.05
24	15.84	4043	v	В	0.00404	1471.29
25	16.08	3797	Endosulfan I	В	0.00130	1259.47
26	16.83	3619		В	0.00362	1123.60
27	17.78	4999		В	0.00500	1256.00
28	18.43	55142		В	0.05514	16855.81
29	19.92	12623		В	0.01262	3570.46
30	20.14	4988	Endrin ketone	В	0.00139	1580.77
31	21.07	23162		В	0.02316	981.49
32	21.18	10373		V	0.01037	2545.22
33	22.32	425185	Decachlorobiphenyl	В	0.01837	99909.18
		1075497			0.40589	298398.09

2002

Software Version	: 6.2.1.0.104:0104 : buf2048: 98599	Date	: 09/25/2009 13:14:07
Operator Sample Number	: tchrom	Sample Name	: 9I20007-BLK1
AutoSampler	BUILT-IN	Rack/Vial	: 1/93 · B
Instrument Serial #	: CN10520010	A/D mV Range	: 1000
Sampling Rate	: 5.000 pts/s	End lime	: 29.45 min
Sample Volume Sample Amount	: 1.000000 ul : 1.0000	Area Reject Dilution Factor	: 1500.000000 : 1.00
Data Acquisition Time	: 09/25/2009 12:09:51	Cycle	: 1

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46193.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46193.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46193.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46193.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46193.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]	
2	8.30	2376		В	0.00238	614.65	SFR
3	8.45	64572		В	0.06457	21865.47	0-1
4	8.53	1541		Е	0.00154	597.91	IV
5	8.80	2583		В	0.00258	789.74	•
6	8.99	5181		В	0.00518	1386.18	
8	9.58	18860		V	0.01886	6258.71	
10	10.71	93550	Tetrachloro-m-xylene (В	0.01071	28741.08	
11	11.28	2247		В	0.00225	608.87	
12	11.53	2948		В	0.00295	684.51	
13	11.72	2631		В	0.00263	677.84	
14	12.17	8234		В	0.00823	2541.92	
17	14.32	6355		В	0.00636	1058.53	
18	14.37	4278	delta-BH C (2C)	V	0.00309	1245.67	
19	15.11	6344		В	0.00634	1992.57	
20	15.53	17376		В	0.01738	5235.06	
	15 62	1611		- 17	0.00464	15/2 /1	

^{247/275} Page 2 of 2

09/25/2009 13:14:07 Result: H:\TURBO6\6890-06\6-SEQ46\6b46193.rst

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
23	16.75	3477	Hept: epoxide (2C)	в	0.00116	659.85
25	18.71	1511		В	0.00151	511.79
26	19.71	3723		В	0.00372	1181.71
27	20.09	19279	Endrin aldehyde (2C)	В	0.00279	5697.46
28	21.16	2967	,	В	0.00297	957.84
31	25.54	172137	Decachlorobiphenyl (2C	В	0.02228	28385.86
		446811			0.19412	113236.63
248/275

Form 1

LCS

ORGANIC ANALYSIS DATA SHEET

8081A

Laboratory:	TestAmerica Bi	uffalo			SDG:			
Client:	Olin Chlor Alka	ali Products - Cle	eveland, TN		Project:	Olin - Charles Gib	son site - NY3A90	25AE03759
Matrix:	<u>Solid</u>	La	boratory ID:	<u>9120007</u>	<u>-BS1</u>	File ID:	<u>6A46181</u>	
Sampled:		Pre	epared:	<u>09/21/09</u>	9 08:00	Analyzed:	<u>09/25/09 04:58</u>	
Solids:		Pre	eparation:	<u>3550B (</u>	<u>GC</u>	Initial/Final:	<u>30.56 g / 10 mL</u>	
Batch:	<u>9120007</u>	Sequence:	<u>RI92924</u>		Calibration:	<u>R9H1803</u>	Instrument:	<u>HP6890-6</u>
CAS NO.	COMPOUND				DILUTION	CONC	. (ug/kg)	Q
319-84-6	alpha-BHC				1	1	4.9	
319-85-7	beta-BHC				1	1	6.7	
319-86-8	delta-BHC				1	1	6.0	
58-89-9	gamma-BHC (L	indane)			1	1	5.6	
SYSTEM MONI	TORING COMP	OUND	ADDED) (ug/kg)	CONC (ug/kg)	% REC	QC LIMITS	Q
Decachlorobipher	nyl		6.:	54	6.38	97	42 - 146	
Tetrachloro-m-xy	lene		6.	54	7.06	108	37 - 135	

* Values outside of QC limits

Software Version	: 6.2.1.0.104:0104	Date	: 09/25/2009 08:17:57
Reprocess Number Operator Sample Number AutoSampler	: buf2048: 98518 : tchrom : : BUILT-IN : HP6800-06	Sample Name Study Rack/Vial Channel	: 9I20007-BS : : 1/81 : A
Instrument Name Instrument Serial # Delay Time	: CN10520010 : 0.00 min	A/D mV Range End Time	: 1000 : 29.50 min
Sample Volume Sample Amount Data Acquisition Time	: 1.00000 ul : 1.0000 : 09/25/2009 04:58:51	Area Reject Dilution Factor Cycle	: 1500.000000 : 1.00 : 6

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6A46181.raw <Modified> Result File : H:\TURBO6\6890-06\6-SEQ46\6A46181.rst

Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6A46181.raw Proc Method : h:\turbo6\6890-06\6a-process.mth from H:\TURBO6\6890-06\6-SEQ46\6A46181.rst Calib Method : h:\turbo6\6890-06\6a-pest(07-23-09)surr(07-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6A46181.rst Report Format File: h:\turbo6\6890-06\6samp.rpt Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP I/ "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	7.24	4339		В	0.00434	1082.01
2	7.84	6154		В	0.00615	2185.58
3	7.92	70710		V	0.07071	24156.72
4	8.06	6093		В	0.00609	1629.16
5	8 38	2270		B	0.00227	591.05
ő	8 73	15027		В	0.01503	4103.41
7	8 97	19896		В	0.01990	7418.44
, 8	9.16	9893		В	0.00989	2446.90
ğ	9.36	342220	Tetrachloro-m-xvlene	В	0.02158	111066.52
10	10.94	1333065	alpha-BHC	В	0.04567	437566.53
11	11 35	12943	alpha anna	В	0.01294	3119.49
12	11 67	5040		В	0.00504	1577.02
13	11 79	1285567	gamma-BHC	v	0.04778	413693.91
14	12.03	566441	beta-BHC	B	0.05113	174032.74
15	12 13	13074		Ē	0.01307	4836.02
16	12.48	1347511	delta-BHC	B	0.04900	427324.85

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09/25/2009 08:17:57 Result: H:\TURBO6\6890-06\6-SEQ46\6A46181.rst

Peak #	Time [mín]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
17	12.60	35968		v	0.03597	9608.93
18	13.00	1282449	Heptachlor	В	0.05038	406744.51
19	13.37	5375		В	0.00538	1521.16
20	13.50	4626		В	0.00463	1607.14
21	13.73	1199517	Aldrin	В	0.04953	377844.12
22	14.03	1508		В	0.00151	605.74
23	14.52	12994		В	0.01299	3021.15
24	14.71	12862		В	0.01286	3966.66
25	14.82	5201		V	0.00520	1361.74
26	14.98	6681		В	0.00668	1655.22
27	15.18	1263871	Hept. epoxide	В	0.05108	378970.70
28	15.46	1280001	gamma chlordane	В	0.04887	386552.55
29	15.76	1122122	alpha chlordane	В	0.04670	342145.89
30	15.96	1247441	4,4'-DDE	В	0.04919	400961.34
31	16.07	939899	Endosulfan I	V	0.03981	276993.92
32	16.61	1274684	Dieldrin	В	0.04690	372991.09
33	16.91	28815		В	0.02881	9122.37
34	17.13	1136168	Endrin	В	0.04664	328215.32
35	17.28	963026	4,4'-DDD	V	0.04457	299431.69
36	17.62	920374	Endosulfan II	В	0.04007	260142.25
37	17.89	989083	4,4'-DDT	В	0.04392	292982.73
38	18.43	56303	,	В	0.05630	17608.58
39	18.54	680201	Endrin aldehyde	V	0.03933	185455.68
40	18.94	480631	Methoxychlor	В	0.04109	143749.23
41	19.49	961367	Endo. Sulfate	В	0.04547	238990.94
42	19.77	4628		В	0.00463	1352.41
43	19.92	13074		В	0.01307	3802.13
44	20.09	1009169	Endrin ketone	В	0.04131	263675.06
45	21.18	13144		В	0.01314	4019.03
46	22.31	450308	Decachlorobiphenyl	В	0.01949	102649.92
		22441732			1.32612	6.73e+06

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98519	Date	: 09/25/2009 08:18:00
Sample Number AutoSampler Instrument Name Instrument Serial # Delay Time Sampling Rate	: tchrom : : BUILT-IN : HP6890-06 : CN10520010 : 0.00 min : 5 0000 pts/s	Sample Name Study Rack/Vial Channel A/D mV Range End Time	: 9i20007-BS : : 1/81 : B : 1000 : 29.50 min
Sample Volume Sample Amount Data Acquisition Time	: 1.000000 ul : 1.0000 : 09/25/2009 04:58:51	Area Reject Dilution Factor Cycle	: 1500.000000 : 1.00 : 6

Raw Data File : H:\TURBO6\6890-06\6-SEQ46\6b46181.raw <Modified>

Result File : H:\TURBO6\6890-06\6-SEQ46\6b46181.rst Inst Method : h:\turbo6\6890-06\6890-6 12-5-08-ins from H:\TURBO6\6890-06\6-SEQ46\6b46181.raw Proc Method : h:\turbo6\6890-06\6b-process.mth from H:\TURBO6\6890-06\6-SEQ46\6b46181.rst Calib Method : h:\turbo6\6890-06\6b-pest(07-23-09)surr(7-23-09)5.mth from H:\TURBO6\6890-06\6-SEQ46\6b46181.rst Sequence File : H:\TURBO6\6890-06\6-SEQ46\6D-46.seq



HP 6890-06 "A" RTXCLP // "B" RTXCLP II

Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1	8.20	1981		B	0.00400	
- 2	8.45	44411		ň	0.00198	724.88
- 4	9.58	12298		В	0.04441	15246.71
5	9.93	4028		В	0.01230	4117.86
6	10.15	3341		В	0.00403	1167.41
7	10.71	132991	Tetrachloro, m vulena (В	0.00334	1035.45
8	11.28	2237	(В	0.01470	41034.68
9	11.52	3199		В	0.00224	577.02
10	12.17	5056		В	0.00320	795.97
11	12.34	7651		В	0.00506	1821.03
12	12.48	561689	alpha-BHC (2C)	в	0.00765	1133.26
13	13.16	2195		Ň	0.03991	176330.26
<u>14</u>	<u>13.44</u>	534734	gamma-BHC (2C)	Б	0.00219	729.21

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09/25/2009 08:18:00 Result: H:\TURBO6\6890-06\6-SEQ46\6b46181.rst

Peak Time Area Component Name BL NG Height (UV) 18 14.59 542674 Heptachlor (2C) B 0.04726 164283.05 20 15.40 473689 Aldrin (2C) B 0.04569 141198.64 21 15.52 21795 V 0.02179 6876.58 23 16.35 3578 B 0.00426 698.00 24 16.75 477451 Hept. epoxide (2C) B 0.04955 134569.45 26 17.19 446156 gamma chlordane (2C) B 0.04955 126523.87 29 17.87 435146 4.4'-DDE (2C) B 0.04995 126523.87 29 17.87 435146 4.4'-DDE (2C) B 0.05001 131938.89 30 18.28 447424 Dieldrin (2C) B 0.05229 101570.36 31 9.15 350604 4.4'-DDT (2C) B 0.04340 87033.88 31 <	. .	_				0 - 0 10 10 - 10 1.15	Į –
1814.59542674Heptachlor (2C)B 0.04726 164283.051915.119907B 0.00991 3131.782015.40473689Aldrin (2C)B 0.00991 3131.782115.5221795V 0.02179 6876.582216.134259B 0.00426 698.002416.75477451Hept. epoxide (2C)B 0.00495 134569.452617.19446156gamma chlordane (2C)B 0.04995 134569.452617.19446156gamma chlordane (2C)B 0.04925 126523.872917.874351464,4'-DDE (2C)B 0.05001 131938.892018.28447424Dieldrin (2C)B 0.04995 12528.063119.153506044,4'-DDD (2C)V 0.04998 101855.813519.853173144,4'-DDT (2C)B 0.04997 65552.503820.15276910Endrin aldehyde (2C)B 0.04989 91646.523720.76319625Endo. Sulfate (2C)B 0.04989 91646.523921.29202100Methoxychlor (2C)V 0.05958 50427.664122.953440B 0.00336 1033.584223.781674B 0.002210 28353.398508089850808914240024353.39	Peak #	Time [min]	Area [uV-sec]	Component Name	BL	NG CONCENTRATION	Height [µV]
1.1.340.3 2.406±06	19 20 22 22 22 22 22 22 22 22 22 22 22 22	14.59 15.11 15.40 15.52 16.13 16.75 17.19 17.53 17.66 17.87 18.28 18.97 19.42 19.85 20.15 20.76 1.15 2.01 2.91 2.91 2.91 3.78 5.53	542674 9907 473689 21795 4259 3578 477451 446156 446223 337633 435146 447424 387091 350604 447424 328088 17314 276910 8319625 8359 202100 N 366253 8440 1674 170837 D 508089	Heptachlor (2C) Aldrin (2C) Hept. epoxide (2C) gamma chlordane (2C) alpha chlordane (2C) Endosulfan I (2C) 4,4'-DDE (2C) Endrin (2C) Endrin (2C) Endrin (2C) Endrin aldehyde (2C) Endrin aldehyde (2C) Endrin aldehyde (2C) Endrin ketone (2C) Methoxychlor (2C) Endrin ketone (2C)	888~8888~888~8888~8888~888~888~888~888	0.04726 0.00991 0.04569 0.02179 0.00426 0.00358 0.04995 0.04715 0.04925 0.03999 0.05001 0.04899 0.05229 0.04998 0.04989 0.04989 0.04989 0.04989 0.04977 0.04512 0.00336 0.05958 0.04530 0.00344 0.00167 0.02210 1.13403	164283.05 3131.78 141198.64 6876.58 698.00 1133.26 134569.45 128229.64 126523.87 94180.48 131938.89 122528.06 101570.36 101855.81 87033.88 91646.52 65552.50 83653.22 1033.58 50427.66 81100.68 893.22 88.90 28353.39

Page 2 of 2

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HOLDING TIME SUMMARY

8081A

Laboratory:	TestAmerica Buff	alo			SDG:	<u>D</u>]	DRAFT						
Client:	<u>Olin Chlor Alkali</u>	Products - Cl	eveland, TN		Project:	<u>0</u>]	in - Charles	<u>Gibson site -</u>	NY3A9025	<u>AE03759</u>			
Sample Name		Date Collected	Date Received	Date Prepared	Days to Prep	Max Days to Prep	Date Analyzed	Days to Analysis	Max Days to Analysis	Q			
CGS-US1-091709)	09/17/09 11:30	09/17/09 15:00	09/21/09 08:00	4	10	09/29/09 02:32	8	40				
CGS-USD1-0917	09	09/17/09 11:40	09/17/09 15:00	09/21/09 08:00	4	10	09/29/09 03:08	8	40				
CGS-DS1-091709)	09/17/09 12:00	09/17/09 15:00	09/21/09 08:00	4	10	09/29/09 03:44	8	40				
CGS-DSD1-0917	09	09/17/09 12:20	09/17/09 15:00	09/21/09 08:00	4	10	09/29/09 04:20	8	40				

					TestAme	erica Buttale	0			111 1/21/2009 2:21:401 M
Matrix: Solid				Prepare	ed using: E	xtractions - 🤅	3550B GC	T \		Surrogate used: 9090238
			Initial	Final		01	ul Sailea	ul	Bar Code	Extraction Comments
Lab Number	Analysis	Prepared	(g)	(mr)	Spike ID	Source ID	opike	Sullogate	Dai Couc	
9120007-BLK1	QC	09/21/09 08:00	30.2	10			0	1000		
9120007-BS1	qc	00:80 60/12/60	30.56	10	9080196		1000	1000		
9120007-MS1	qc	09/21/09 08:00	30.2	10	9080196	RS10643-03	0001	1000		
9120007-MSD1	QC	09/21/09 08:00	30.04	10	9080196	RSI0643-03	0001	1000		
RS10482-01	8081	09/21/09 08:00	30.42	10				1000		
RSI0634-01	8081_ASP00	09/21/09 08:00	30.28	10				1000		
RSI0634-02	8081_ASP00	00/21/09 08:00	30.15	10				1000		
RSI0634-03	8081_ASP00	06/21/09 08:00	30.54	10				1000		
RSI0634-04	8081_ASP00	09/21/09 08:00	30.44	10				1000	NU INNI INNI AVANA AVANA ILANA ILANA	
RSI0643-01	8081	09/21/09 08:00	30.48	10				1000	NATIONAL AND	
RSI0643-02	8081	00:80 60/12/60	30.02	10				1000		
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Chr Cleanup: Elmal Concentration: Vialed: Concentration: Lentered: CXM Closed: CM ξ surrogated: Spiked: Extracted:

Spiking Witnessed By

Date

Preparation Reviewed By

Date

Extracts Received By

Date

Page l of l

9120007

PREPARATION BENCH SHEET

'rinted: 9/21/2009 2:21:40PM

		Sequence 42	Comments												40K Baldehud. L	AOK BOK									Date:
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		rument ID. vook #	e & Initial	23-09	9413																			Rev.0 12/20/	

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Software Version	: 6.2.1.0.104:0104	Date	: 09/24/2009 12:44:46
Reprocess Number	: buf2048: 98450		
Operator	: tchrom	Sample Name	: RSI0634-01
Sample Number	:	Study	:
AutoSampler	: NONE	Rack/Vial	: 0/0
Instrument Name	: HP5890-16	Channel	: B
Interface Serial #	: 3090270361	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 7.00 min
Sampling Rate	: 16.6660 pts/s		
Sample Volume	: 1.000000 uL	Area Reject	: 1000.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/24/2009 09:26:40	Cycle	: 1

Raw Data File : H:\TURBO6\5890-16\16SEQ35\16b35030.raw <Modified>

Result File : H:\TURBO6\5890-16\16SEQ35\16b35030.rst

Inst Method : H:\TURBO6\5890-16\PPCBINS from H:\TURBO6\5890-16\16SEQ35\16b35030.raw Proc Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35030.rst Calib Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35030.rst Report Format File: h:\turbo6\5890-16\16screen.rpt Sequence File : H:\TURBO6\5890-16\16SEQ35\16D35.seq



PCB Screen

09/24/2009 12:44:46 Result: H:\TURBO6\5890-16\16SEQ35\16b35030.rst

Pe	eak #	Ret Time [min]	BL	Dilution Needed	Area [uV-sec]	Component Name	NG conc.	Dilution Needed
		1.30		0	87953	AR1242	0.00731	0
		1.60		0	93627	AR1248	0.00546	0
		2.10		0	5674	AR1254	3.47e-04	0
					187254			

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98452	Date	: 09/24/2009 12:44:50
Operator	: tchrom	Sample Name	: RSI0634-02
Sample Number	:	Study	:
AutoSampler	: NONE	Rack/Vial	: 0/0
Instrument Name	: HP5890-16	Channel	: B
Interface Serial #	: 3090270361	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 7.00 min
Sampling Rate	: 16.6660 pts/s		
Sample Volume	: 1.000000 uL	Area Reject	: 1000.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/24/2009 09:37:38	Cycle	: 2

Raw Data File : H:\TURBO6\5890-16\16SEQ35\16b35031.raw <Modified>

Result File : H:\TURBO6\5890-16\16SEQ35\16b35031.rst

Inst Method : H:\TURBO6\5890-16\PPCBINS from H:\TURBO6\5890-16\16SEQ35\16b35031.raw Proc Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35031.rst Calib Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35031.rst Report Format File: h:\turbo6\5890-16\16screen.rpt Sequence File : H:\TURBO6\5890-16\16SEQ35\16D35.seq



PCB Screen

09/24/2009 12:44:50 Result: H:\TURBO6\5890-16\16SEQ35\16b35031.rst

Peak #	Ret Time [min]	BL	Dilution Needed	Area [uV-sec]	Component Name	NG conc.	Dilution Needed
 	1.30		0	408361	AR1242	0.03392	0
	1.60		0	408361	AR1248	0.02383	0
				816723			

Software Version Reprocess Number	: 6.2.1.0.104:0104 : buf2048: 98454	Date	: 09/24/2009 12:44:54
Operator	: tchrom	Sample Name	: RSI0634-03
Sample Number	:	Study	:
AutoSampler	: NONE	Rack/Vial	: 0/0
Instrument Name	: HP5890-16	Channel	: B
Interface Serial #	: 3090270361	A/D mV Range	: 1000
Delay Time	: 0.00 min	End Time	: 7.00 min
Sampling Rate	: 16.6660 pts/s		
Sample Volume	: 1.000000 uL	Area Reject	: 1000.000000
Sample Amount	: 1.0000	Dilution Factor	: 10.00
Data Acquisition Time	: 09/24/2009 09:48:32	Cycle	: 3

Raw Data File : H:\TURBO6\5890-16\16SEQ35\16b35032.raw <Modified>

Result File : H:\TURBO6\5890-16\16SEQ35\16b35032.rst

Inst Method : H:\TURBO6\5890-16\PPCBINS from H:\TURBO6\5890-16\16SEQ35\16b35032.raw Proc Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35032.rst Calib Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35032.rst Report Format File: h:\turbo6\5890-16\16screen.rpt Sequence File : H:\TURBO6\5890-16\16SEQ35\16D35.seq



PCB Screen

09/24/2009 12:44:54 Result: H:\TURBO6\5890-16\16SEQ35\16b35032.rst

Peak #	Ret Time [min]	BL	Dilution Needed	Area [uV-sec]	Component Name	NG conc.	Dilution Needed
 	1.30		0	84216	AR1242	0.00699	0
	1.60		0	84216	AR1248	0.00491	0
				168431			

7		
: 6.2.1.0.104:0104	Date	: 09/24/2009 12:44:58
: buf2048: 98456		
: tchrom	Sample Name	: RSI0634-04
:	Study	:
: NONE	Rack/Vial	: 0/0
: HP5890-16	Channel	: B
: 3090270361	A/D mV Range	: 1000
: 0.00 min	End Time	: 7.00 min
: 16.6660 pts/s		
: 1.000000 uL	Area Reject	: 1000.000000
: 1.0000	Dilution Factor	: 10.00
: 09/24/2009 09:59:51	Cycle	: 4
	: 6.2.1.0.104:0104 : buf2048: 98456 : tchrom : : NONE : HP5890-16 : 3090270361 : 0.00 min : 16.6660 pts/s : 1.00000 uL : 1.0000 : 09/24/2009 09:59:51	: 6.2.1.0.104:0104 Date : buf2048: 98456 : tchrom Sample Name : Study : NONE Rack/Vial : HP5890-16 Channel : 3090270361 A/D mV Range : 0.00 min End Time : 16.6660 pts/s : 1.00000 uL Area Reject : 1.0000 Dilution Factor : 09/24/2009 09:59:51 Cycle

Raw Data File : H:\TURBO6\5890-16\16SEQ35\16b35033.raw <Modified>

Result File : H:\TURBO6\5890-16\16SEQ35\16b35033.rst

Inst Method : H:\TURBO6\5890-16\PPCBINS from H:\TURBO6\5890-16\16SEQ35\16b35033.raw Proc Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35033.rst Calib Method : h:\turbo6\5890-16\16-b-screen.mth from H:\TURBO6\5890-16\16SEQ35\16b35033.rst Report Format File: h:\turbo6\5890-16\16screen.rpt Sequence File : H:\TURBO6\5890-16\16SEQ35\16D35.seq



PCB Screen

09/24/2009 12:44:58 Result: H:\TURBO6\5890-16\16SEQ35\16b35033.rst

Peak #	Ret Time [min]	BL	Dilution Needed	Area [uV-sec]	Component Name	NG conc.	Dilution Needed
 	1.30		0	76676	AR1242	0.00637	0
	1.60		0	76676	AR1248	0.00447	0

153353

TestAmerica Buffalo SDG: CLASS: WET_BATCH METHOD: Dry Weight

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ANALYSES DATA PACKAGE COVER PAGE

Laboratory:	TestAmerica Buffalo	SDG:
Client:	Olin Chlor Alkali Products - Cleveland, TN	Project: Olin - Charles Gibson site - NY3A9025AE03759
	Client Sample Id:	Lab Sample Id:
	CGS-US1-091709	<u>RSI0634-01</u>
	CGS-USD1-091709	<u>RSI0634-02</u>
	CGS-DS1-091709	<u>RSI0634-03</u>
	CGS-DSD1-091709	<u>RSI0634-04</u>

METHOD DETECTION AND REPORTING LIMITS

Laboratory:	TestAmerica Buffalo			SDG:	
Client:	Olin Chlor Alkali Products - Clevela	<u>nd, TN</u>	P	roject: <u>Olin - Ch</u>	arles Gibson site - NY3A9025AE(
Matrix:	Solid		Instr	ument: <u>INST</u>	
	Analyte	MDL	MRL	Units	
	Percent Solids	0.010	0.010	%	

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CGS-US1-091709

INORGANIC ANALYSIS DATA SHEET

Laboratory:	<u>TestAmerica Bu</u>	ffalo		SDG:	SDG:				
Client:	Olin Chlor Alkal	i Products - Cleveland, T	N	Project: (Project: Olin - Charles Gibson site - NY3A9025AE03759				
Matrix:	<u>Solid</u>	Laborato	ry ID:	<u>RSI0634-01</u>	File ID:				
Sampled:	<u>09/17/09 11:30</u>	Pre	Prepared: <u>09/18/09 09:20</u> Analyzed:				: <u>09/18/09 15:15</u>		
Solids:	<u>26.93</u>	Prepar	Preparation: Dry Weight Initial/Final				l: <u>10 g / 10 g</u>		
Batch:	<u>9I18022</u>	Sequence:			Calibration:	Instrur	Instrument: INST		
CAS NO.	Analyte		Concentration		Units	Dilution Factor	Q	Method	
STL00234	Percent Solids			27	%	% 1		Dry Weight	

CGS-USD1-091709

INORGANIC ANALYSIS DATA SHEET

Laboratory:	<u>TestAmerica Bu</u>	<u>ffalo</u>			SDG:					
Client:	Olin Chlor Alkal	li Products - C	Cleveland, T	N	Project: Olin - Charles Gibson site - NY3A9025AE03759					
Matrix:	<u>Solid</u>		Laborato	ry ID:	<u>RSI0634-02</u>		File ID:			
Sampled:	<u>09/17/09 11:40</u>		Рге	pared:	<u>09/18/09 09:20</u>		Analyzed:	: <u>09/18/09 15:17</u>		
Solids:	<u>27.37</u>		Prepar	ation:	Dry Weight		Initial/Final:	<u>10 g / 10 g</u>	L	
Batch:	<u>9118022</u>	Sequence:			(Calibration:		Instrur	nent: <u>INST</u>	
CAS NO.	Analyte			С	oncentration	Units	Dilution Factor	Q	Method	
STL00234	Percent Solids				27	%	1		Dry Weight	

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CGS-DS1-091709

INORGANIC ANALYSIS DATA SHEET

Dry Weight

Laboratory:	<u>TestAmerica Bu</u>	<u>ffalo</u>				SDG:				
Client:	Olin Chlor Alka	li Products - (<u>Cleveland, T</u>	N	Project: Olin - Charles Gibson site - NY3A9025AE03759					
Matrix:	Solid		Laborato	ry ID:	<u>RSI0634-03</u>		File ID:			
Sampled:	09/17/09 12:00		Prej	pared:	<u>09/18/09 09:20</u>	1	Analyzed:	<u>09/18/09 1</u>	<u>.5:19</u>	
Solids:	27.16		Prepar	ation:	Dry Weight		Initial/Final:	<u>10 g / 10 g</u>	L.	
Batch:	<u>9118022</u>	Sequence:				Calibration:		Instrun	nent: INST	
CAS NO.	Analyte			С	Concentration	Units	Dilution Factor	Q	Method	
STL00234	Percent Solids				27	%	1		Dry Weight	

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CGS-DSD1-091709

INORGANIC ANALYSIS DATA SHEET

Laboratory:	TestAmerica Bu	ffalo		SDG:					
Client:	<u>Olin Chlor Alka</u>	li Products - Cleveland, 7	N	Project: Olin - Charles Gibson site - NY3A9025AE03759					
Matrix:	<u>Solid</u>	Laborato	ry ID:	<u>RSI0634-04</u>		File ID:			
Sampled:	<u>09/17/09 12:20</u>	Pre	pared:	<u>09/18/09 09:20</u>		Analyzed:	: <u>09/18/09 15:21</u>		
Solids:	<u>37.92</u>	Ргера	ration:	Dry Weight		Initial/Final:	<u>10 g / 10 g</u>	S	
Batch:	<u>9I18022</u>	Sequence:			Calibration:		Instrum	nent: <u>INST</u>	
CAS NO.	Analyte		c	Concentration	Units	Dilution Factor	Q	Method	
STL00234	Percent Solids			38	%	% 1		Dry Weight	

PREPARATION BATCH SUMMARY

<u>i estAmerica i</u>	surraio			SDG:							
Client: <u>Olin Chlor Alkali Product</u>			<u>is - Cleveland, TN</u> Project:					Olin - Charles Gibson site - NY3A9025AE03759			
h: <u>9118022</u> Batch Matrix: <u>Solid</u>					Preparation:			Dry Weight			
AMPLE NAME	I	LAB S	AMPLE ID	INIT	IAL	FINA	L	DATE PREPARED	TOT/DIS		
GS-US1-091709		RSI	0634-01	10.0	g	10.0	g	09/18/09 09:20	Total		
GS-USD1-091709		RSI	0634-02	10.0	g	10.0	g	09/18/09 09:20	Total		
GS-DS1-091709		RSI	0634-03	10.0	g	10.0	g	09/18/09 09:20	Total		
GS-DSD1-091709		RSI	0634-04	10.0	g	10.0	g	09/18/09 09:20	Total		
	<u>Olin Chlor Alk</u> <u>9118022</u> MPLE NAME GS-US1-091709 GS-USD1-091709 GS-DS1-091709	Olin Chlor Alkali Products - (9118022 Batch Matr IMPLE NAME I GS-US1-091709 GS-DS1-091709 GS-DSD1-091709 GS-DSD1-091709	Olin Chlor Alkali Products - Clevels9118022Batch Matrix:MPLE NAMELAB SGS-US1-091709RSHGS-USD1-091709RSHGS-DS1-091709RSHGS-DSD1-091709RSH	Itervinence partatoQlin Chlor Alkali Products - Cleveland, TN9118022Batch Matrix:SolidMPLE NAMELAB SAMPLE IDGS-US1-091709RSI0634-01GS-USD1-091709RSI0634-02GS-DS1-091709RSI0634-03GS-DSD1-091709RSI0634-04	Olin Chlor Alkali Products - Cleveland, TN 9118022 Batch Matrix: Solid MPLE NAME LAB SAMPLE ID INIT. GS-US1-091709 RSI0634-01 10.0 GS-DS1-091709 RSI0634-03 10.0 GS-DSD1-091709 RSI0634-04 10.0	Olin Chlor Alkali Products - Cleveland, TNProject9118022Batch Matrix:SolidMPLE NAMELAB SAMPLE IDINITIALGS-US1-091709RSI0634-0110.0 gGS-USD1-091709RSI0634-0210.0 gGS-DSD1-091709RSI0634-0310.0 g	Olin Chlor Alkali Products - Cleveland, TN Project: 9118022 Batch Matrix: Solid Preparation: MPLE NAME LAB SAMPLE ID INITIAL FINA GS-US1-091709 RSI0634-01 10.0 g 10.0 GS-USD1-091709 RSI0634-02 10.0 g 10.0 GS-DSD1-091709 RSI0634-03 10.0 g 10.0	Olin Chlor Alkali Products - Cleveland, TNProject:Olin9118022Batch Matrix:SolidPreparation:DryMPLE NAMELAB SAMPLE IDINITIALFINALGS-US1-091709RSI0634-0110.0g10.0gGS-USD1-091709RSI0634-0210.0g10.0gGS-DSD1-091709RSI0634-0310.0g10.0g	Olin Chlor Alkali Products - Cleveland, TN Project: Olin - Charles Gibson site - NY 9118022 Batch Matrix: Solid Preparation: Dry Weight MPLE NAME LAB SAMPLE ID INITIAL FINAL DATE PREPARED GS-US1-091709 RSI0634-01 10.0 g 10.0 g 09/18/09 09:20 GS-USD1-091709 RSI0634-02 10.0 g 10.0 g 09/18/09 09:20 GS-DS1-091709 RSI0634-03 10.0 g 10.0 g 09/18/09 09:20 GS-DSD1-091709 RSI0634-04 10.0 g 10.0 g 09/18/09 09:20		