

Cover Page

Order ID : Q2458

Project ID : South River WM Replacement

Client : CDM Smith

Lab Sample Number

Q2458-01
Q2458-02
Q2458-03
Q2458-04
Q2458-05
Q2458-06
Q2458-07
Q2458-08
Q2458-09
Q2458-10

Client Sample Number

TP-76
TP-55
TP-68
TP-67
TP-66
TP-60
TP-62
TP-63
TP-59
FB-06272025

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hard copy data package has been authorized by the laboratory manager or his designee, as verified by the following signature.

Signature : _____

Date: 7/9/2025



284 Sheffield Street, Mountainside, NJ 07092
Phone: 908 789 8900 Fax: 908 789 8922

CASE NARRATIVE

CDM Smith

Project Name: South River WM Replacement

Project # N/A

Order ID # Q2458

Test Name: Pesticide-TCL

A. Number of Samples and Date of Receipt:

9 Solid samples were received on 06/27/2025.

1 Water sample was received on 06/27/2025.

B. Parameters

According to the Chain of Custody document, the following analyses were requested: Diesel Range Organics, Gasoline Range Organics, Herbicide, Mercury, Metals ICP-TAL, METALS-TAL, PCB, Pesticide-TCL, SVOC-TCL BNA -20, VOC-TCLVOA-10 and VOC-TCLVOA-10. This data package contains results for Pesticide-TCL.

C. Analytical Techniques:

The analysis was performed on instrument ECD_D. The front column is ZB-MR1 which is 30 meters, 0.32 mm ID, 0.5 um df.; Catalog # 7HM-G016-17. The rear column is ZB-MR2 which is 30 meters, 0.32 mm ID, 0.25 um df, Catalog #: 7HMG017- 11. The analysis of Pesticide-TCLs was based on method 8081B and extraction was done based on method 3510.

D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria.

The Retention Times were acceptable for all samples.

The MS recoveries met the requirements for all compounds due to matrix interference.

The MSD recoveries met the acceptable requirements due to matrix interference.

The RPD met criteria .

The Blank Spike met requirements for all samples .

The Blank Spike Duplicate met requirements for all samples .

The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements .

The Continuous Calibration met the requirements .

E. Additional Comments:

Sample TP-60 was reported with J flag on form 1 for compound 4,4-DDD based on reporting criteria of high concentration from both column. Now for other column compound detection is below MDL therefore it is not detecting on form 10.



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Sample TP-62 was reported with J flag on form 1 for compound Heptachlor based on reporting criteria of high concentration from both column. Now for other column compound detection is below MDL therefore it is not detecting on form 10. The soil samples results are based on a dry weight basis.

F. Manual Integration Comments:

Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. The laboratory manager or his designee, as verified by the following signature has authorized release of the data contained in this hard copy data package.

Signature_____

DATA REPORTING QUALIFIERS- ORGANIC

For reporting results, the following “ Results Qualifiers” are used:

Value	If the result is a value greater than or equal to the detection limit, report the value
U	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. “10 U”. This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
ND	Indicates the analyte was analyzed for, but not detected
J	Indicates an estimated value. This flag is used: (1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.) (2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3 ug/L was calculated report as 3 J. This flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others.
B	Indicates the analyte was found in the blank as well as the sample report as “12 B”.
E	Indicates the analyte ‘s concentration exceeds the calibrated range of the instrument for that specific analysis.
D	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
P	This flag is used for Pesticide/PCB target analyte when there is >25% difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form 1 and flagged with a “P”.
N	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
A	This flag indicates that a Tentatively Identified Compound is a suspected aldol-condensation product.
Q	Indicates the LCS did not meet the control limits requirements

APPENDIX A

QA REVIEW GENERAL DOCUMENTATION

Project #: Q2458

Completed

For thorough review, the report must have the following:

GENERAL:

Are all original paperwork present (chain of custody, record of communication,airbill, sample management lab chronicle, login page)

✓

Check chain-of-custody for proper relinquish/return of samples

✓

Is the chain of custody signed and complete

✓

Check internal chain-of-custody for proper relinquish/return of samples /sample extracts

✓

Collect information for each project id from server. Were all requirements followed

✓

COVER PAGE:

Do numbers of samples correspond to the number of samples in the Chain of Custody on login page

✓

Do lab numbers and client Ids on cover page agree with the Chain of Custody

✓

CHAIN OF CUSTODY:

Do requested analyses on Chain of Custody agree with form I results

✓

Do requested analyses on Chain of Custody agree with the log-in page

✓

Were the correct method log-in for analysis according to the Analytical Request and Chain of Castody

✓

Were the samples received within hold time

✓

Were any problems found with the samples at arrival recorded in the Sample Management Laboratory Chronicle

✓

ANALYTICAL:

Was method requirement followed?

✓

Was client requirement followed?

✓

Does the case narrative summarize all QC failure?

✓

All runlogs and manual integration are reviewed for requirements

✓

All manual calculations and /or hand notations verified

✓

QA Review Signature: PRADIP PRAJAPATI

Date: 07/09/2025

LAB CHRONICLE

OrderID:	Q2458	OrderDate:	6/27/2025 4:22:00 PM
Client:	CDM Smith	Project:	South River WM Replacement
Contact:	Marcie Ann Encinas	Location:	D51,VOA Ref. #2 Soil,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q2458-01	TP-76	SOIL			06/26/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-02	TP-55	SOIL			06/26/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Diesel Range Organics	8015D		07/02/25	07/03/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-03	TP-68	SOIL			06/27/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-04	TP-67	SOIL			06/27/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-05	TP-66	SOIL			06/27/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Diesel Range Organics	8015D		07/02/25	07/03/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	

LAB CHRONICLE

Q2458-06	TP-60	SOIL			06/27/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Diesel Range Organics	8015D		07/02/25	07/03/25	
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-07	TP-62	SOIL			06/27/25			06/27/25
			Gasoline Range Organics	8015D			07/01/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-08	TP-63	SOIL			06/27/25			06/27/25
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/01/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-09	TP-59	SOIL			06/27/25			06/27/25
			Gasoline Range Organics	8015D			06/30/25	
			PCB	8082A		07/01/25	07/02/25	
			Pesticide-TCL	8081B		07/01/25	07/01/25	
Q2458-10	FB-06272025	Water			06/27/25			06/27/25
			Diesel Range Organics	8015D		07/02/25	07/02/25	
			Gasoline Range Organics	8015D			07/01/25	
			PCB	8082A		07/02/25	07/02/25	
			Pesticide-TCL	8081B		07/03/25	07/03/25	

Hit Summary Sheet
SW-846

SDG No.: Q2458

Order ID: Q2458

Client: CDM Smith

Project ID: South River WM Replacement

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID : TP-76								
Q2458-01	TP-76	SOIL	Dieldrin	1.90	P	0.15	1.90	ug/kg
Q2458-01	TP-76	SOIL	Endrin	0.66	J	0.15	1.90	ug/kg
Q2458-01	TP-76	SOIL	4,4-DDT	1.20	JP	0.15	1.90	ug/kg
Total Concentration:				3.760				
Client ID : TP-67								
Q2458-04	TP-67	SOIL	4,4-DDE	0.34	JP	0.16	1.90	ug/kg
Q2458-04	TP-67	SOIL	alpha-Chlordane	0.48	JP	0.13	1.90	ug/kg
Q2458-04	TP-67	SOIL	gamma-Chlordane	0.29	J	0.17	1.90	ug/kg
Total Concentration:				1.110				
Client ID : TP-66								
Q2458-05	TP-66	SOIL	4,4-DDT	0.40	JP	0.16	1.90	ug/kg
Total Concentration:				0.400				
Client ID : TP-60								
Q2458-06	TP-60	SOIL	4,4-DDD	0.17	J	0.16	1.80	ug/kg
Q2458-06	TP-60	SOIL	4,4-DDT	0.36	JP	0.15	1.80	ug/kg
Q2458-06	TP-60	SOIL	alpha-Chlordane	0.90	JP	0.13	1.80	ug/kg
Q2458-06	TP-60	SOIL	gamma-Chlordane	0.46	J	0.16	1.80	ug/kg
Total Concentration:				1.890				
Client ID : TP-62								
Q2458-07	TP-62	SOIL	Heptachlor	0.17	J	0.13	1.90	ug/kg
Total Concentration:				0.170				
Client ID : TP-63								
Q2458-08	TP-63	SOIL	4,4-DDD	0.41	J	0.17	2.00	ug/kg
Q2458-08	TP-63	SOIL	4,4-DDT	0.44	JP	0.16	2.00	ug/kg
Q2458-08	TP-63	SOIL	alpha-Chlordane	0.58	J	0.14	2.00	ug/kg
Q2458-08	TP-63	SOIL	gamma-Chlordane	0.40	J	0.17	2.00	ug/kg
Total Concentration:				1.830				



QC SUMMARY

Surrogate Summary

SDG No.: Q2458

Client: CDM Smith

Analytical Method: 8081B

Lab Sample ID	Client ID	Parameter	Column	Spike	Result	Recovery(%)	Qual	Limits(%)	
								Low	High
I.BLK-PD088990.D	PIBLK-PD088990.D	Decachlorobiphen	1	20	18.2	91		57	171
		Tetrachloro-m-xyl	1	20	16.1	80		61	148
		Decachlorobiphen	2	20	18.6	93		57	171
		Tetrachloro-m-xyl	2	20	17.7	88		61	148
I.BLK-PD089264.D	PIBLK-PD089264.D	Decachlorobiphen	1	20	22.9	115		57	171
		Tetrachloro-m-xyl	1	20	20.6	103		61	148
		Decachlorobiphen	2	20	23.3	117		57	171
		Tetrachloro-m-xyl	2	20	22.7	114		61	148
PB168672BL	PB168672BL	Decachlorobiphen	1	20	16.9	85		20	144
		Tetrachloro-m-xyl	1	20	16.9	84		19	148
		Decachlorobiphen	2	20	17.3	87		20	144
		Tetrachloro-m-xyl	2	20	20.0	100		19	148
PB168672BS	PB168672BS	Decachlorobiphen	1	20	20.6	103		20	144
		Tetrachloro-m-xyl	1	20	18.3	92		19	148
		Decachlorobiphen	2	20	19.5	98		20	144
		Tetrachloro-m-xyl	2	20	20.8	104		19	148
I.BLK-PD089276.D	PIBLK-PD089276.D	Decachlorobiphen	1	20	23.0	115		57	171
		Tetrachloro-m-xyl	1	20	20.9	105		61	148
		Decachlorobiphen	2	20	22.4	112		57	171
		Tetrachloro-m-xyl	2	20	23.2	116		61	148
Q2458-01	TP-76	Decachlorobiphen	1	20	13.9	70		20	144
		Tetrachloro-m-xyl	1	20	16.5	83		19	148
		Decachlorobiphen	2	20	13.1	66		20	144
		Tetrachloro-m-xyl	2	20	17.8	89		19	148
Q2458-02	TP-55	Decachlorobiphen	1	20	9.41	47		20	144
		Tetrachloro-m-xyl	1	20	11.0	55		19	148
		Decachlorobiphen	2	20	7.45	37		20	144
		Tetrachloro-m-xyl	2	20	11.8	59		19	148
Q2458-03	TP-68	Decachlorobiphen	1	20	13.1	66		20	144
		Tetrachloro-m-xyl	1	20	16.2	81		19	148
		Decachlorobiphen	2	20	10.9	55		20	144
		Tetrachloro-m-xyl	2	20	17.9	89		19	148
I.BLK-PD089287.D	PIBLK-PD089287.D	Decachlorobiphen	1	20	22.1	110		57	171
		Tetrachloro-m-xyl	1	20	21.0	105		61	148
		Decachlorobiphen	2	20	20.1	100		57	171
		Tetrachloro-m-xyl	2	20	23.4	117		61	148
Q2458-04	TP-67	Decachlorobiphen	1	20	14.9	75		20	144
		Tetrachloro-m-xyl	1	20	19.1	95		19	148
		Decachlorobiphen	2	20	15.0	75		20	144
		Tetrachloro-m-xyl	2	20	20.0	100		19	148
Q2458-04MS	TP-67MS	Decachlorobiphen	1	20	16.3	82		20	144

Surrogate Summary

SDG No.: **Q2458**

Client: **CDM Smith**

Analytical Method: **8081B**

Lab Sample ID	Client ID	Parameter	Column	Spike	Result	Recovery(%)	Qual	Limits(%)	
								Low	High
Q2458-04MS	TP-67MS	Tetrachloro-m-xyl	1	20	19.5	98		19	148
		Decachlorobiphen	2	20	16.5	82		20	144
		Tetrachloro-m-xyl	2	20	21.3	107		19	148
Q2458-04MSD	TP-67MSD	Decachlorobiphen	1	20	16.2	81		20	144
		Tetrachloro-m-xyl	1	20	19.7	98		19	148
		Decachlorobiphen	2	20	16.6	83		20	144
		Tetrachloro-m-xyl	2	20	21.6	108		19	148
Q2458-05	TP-66	Decachlorobiphen	1	20	10.6	53		20	144
		Tetrachloro-m-xyl	1	20	11.6	58		19	148
		Decachlorobiphen	2	20	10.4	52		20	144
		Tetrachloro-m-xyl	2	20	12.6	63		19	148
Q2458-06	TP-60	Decachlorobiphen	1	20	11.6	58		20	144
		Tetrachloro-m-xyl	1	20	12.4	62		19	148
		Decachlorobiphen	2	20	7.43	37		20	144
		Tetrachloro-m-xyl	2	20	13.7	68		19	148
Q2458-07	TP-62	Decachlorobiphen	1	20	12.5	62		20	144
		Tetrachloro-m-xyl	1	20	15.8	79		19	148
		Decachlorobiphen	2	20	10.5	52		20	144
		Tetrachloro-m-xyl	2	20	17.6	88		19	148
Q2458-08	TP-63	Decachlorobiphen	1	20	10.2	51		20	144
		Tetrachloro-m-xyl	1	20	11.6	58		19	148
		Decachlorobiphen	2	20	7.67	38		20	144
		Tetrachloro-m-xyl	2	20	12.7	64		19	148
Q2458-09	TP-59	Decachlorobiphen	1	20	12.9	65		20	144
		Tetrachloro-m-xyl	1	20	17.6	88		19	148
		Decachlorobiphen	2	20	11.4	57		20	144
		Tetrachloro-m-xyl	2	20	19.0	95		19	148
I.BLK-PD089298.D	PIBLK-PD089298.D	Decachlorobiphen	1	20	21.1	106		57	171
		Tetrachloro-m-xyl	1	20	21.6	108		61	148
		Decachlorobiphen	2	20	18.2	91		57	171
		Tetrachloro-m-xyl	2	20	24.2	121		61	148
I.BLK-PD089326.D	PIBLK-PD089326.D	Decachlorobiphen	1	20	19.1	96		57	171
		Tetrachloro-m-xyl	1	20	19.5	97		61	148
		Decachlorobiphen	2	20	18.0	90		57	171
		Tetrachloro-m-xyl	2	20	21.9	109		61	148
PB168718BL	PB168718BL	Decachlorobiphen	1	20	18.8	94		57	171
		Tetrachloro-m-xyl	1	20	18.2	91		61	148
		Decachlorobiphen	2	20	19.0	95		57	171
		Tetrachloro-m-xyl	2	20	20.4	102		61	148
PB168718BS	PB168718BS	Decachlorobiphen	1	20	18.6	93		57	171
		Tetrachloro-m-xyl	1	20	17.7	89		61	148

Surrogate Summary

SDG No.: Q2458

Client: CDM Smith

Analytical Method: 8081B

Lab Sample ID	Client ID	Parameter	Column	Spike	Result	Recovery(%)	Qual	Limits(%)	
								Low	High
PB168718BS	PB168718BS	Decachlorobiphen	2	20	18.9	95		57	171
		Tetrachloro-m-xyl	2	20	20.7	103		61	148
PB168718BSD	PB168718BSD	Decachlorobiphen	1	20	18.4	92		57	171
		Tetrachloro-m-xyl	1	20	17.6	88		61	148
		Decachlorobiphen	2	20	19.0	95		57	171
		Tetrachloro-m-xyl	2	20	20.7	104		61	148
Q2458-10	FB-06272025	Decachlorobiphen	1	20	13.2	66		57	171
		Tetrachloro-m-xyl	1	20	18.5	93		61	148
		Decachlorobiphen	2	20	13.7	69		57	171
		Tetrachloro-m-xyl	2	20	21.0	105		61	148
I.BLK-PD089337.D	PIBLK-PD089337.D	Decachlorobiphen	1	20	19.0	95		57	171
		Tetrachloro-m-xyl	1	20	19.8	99		61	148
		Decachlorobiphen	2	20	19.1	95		57	171
		Tetrachloro-m-xyl	2	20	22.4	112		61	148

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q2458</u>	Analytical Method:	<u>8081B</u>
Client:	<u>CDM Smith</u>	DataFile :	<u>PD089291.D</u>

	Parameter	Spike	Sample		Units	Rec	Rec Qual	RPD	RPD Qual	Low	Limits	
			Result								High	RPD
Lab Sample ID:	Q2458-04MS (Column 1)	Client Sample ID:		TP-67MS								
	alpha-BHC	18.56	0	17.7	ug/kg	95				60	144	
	beta-BHC	18.56	0	17.3	ug/kg	93				54	143	
	delta-BHC	18.56	0	17.9	ug/kg	96				29	151	
	gamma-BHC (Lindane)	18.56	0	17.5	ug/kg	94				61	140	
	Heptachlor	18.56	0	16.5	ug/kg	89				63	135	
	Aldrin	18.56	0	18.0	ug/kg	97				49	139	
	Heptachlor epoxide	18.56	0	17.6	ug/kg	95				41	156	
	Endosulfan I	18.56	0	17.8	ug/kg	96				56	142	
	Dieldrin	18.56	0	17.6	ug/kg	95				47	161	
	4,4'-DDE	18.56	0.2077	18.0	ug/kg	95				55	136	
	Endrin	18.56	0	16.6	ug/kg	89				57	139	
	Endosulfan II	18.56	0	17.1	ug/kg	92				40	163	
	4,4'-DDD	18.56	0	18.4	ug/kg	99				47	163	
	Endosulfan sulfate	18.56	0	16.4	ug/kg	88				62	139	
	4,4'-DDT	18.56	0	14.0	ug/kg	75				51	146	
	Methoxychlor	18.56	0	13.4	ug/kg	72				54	136	
	Endrin ketone	18.56	0	17.2	ug/kg	93				60	129	
	Endrin aldehyde	18.56	0	16.8	ug/kg	91				59	132	
	alpha-Chlordane	18.56	0.4784	18.0	ug/kg	94				39	166	
	gamma-Chlordane	18.56	0.2856	18.2	ug/kg	97				44	175	
Lab Sample ID:	Q2458-04MS (Column 2)	Client Sample ID:		TP-67MS								
	alpha-BHC	18.56	0	17.7	ug/kg	95				60	144	
	beta-BHC	18.56	0	17.6	ug/kg	95				54	143	
	delta-BHC	18.56	0	16.9	ug/kg	91				29	151	
	gamma-BHC (Lindane)	18.56	0	17.5	ug/kg	94				61	140	
	Heptachlor	18.56	0	16.3	ug/kg	88				63	135	
	Aldrin	18.56	0	17.9	ug/kg	96				49	139	
	Heptachlor epoxide	18.56	0	18.2	ug/kg	98				41	156	
	Endosulfan I	18.56	0	17.6	ug/kg	95				56	142	
	Dieldrin	18.56	0	17.7	ug/kg	95				47	161	
	4,4'-DDE	18.56	0.3449	17.9	ug/kg	96				55	136	
	Endrin	18.56	0	16.9	ug/kg	91				57	139	
	Endosulfan II	18.56	0	17.2	ug/kg	93				40	163	
	4,4'-DDD	18.56	0	17.0	ug/kg	92				47	163	
	Endosulfan sulfate	18.56	0	16.1	ug/kg	87				62	139	

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q2458</u>	Analytical Method:	<u>8081B</u>
Client:	<u>CDM Smith</u>	DataFile :	<u>PD089291.D</u>

Parameter	Spike	Sample		Units	Rec	Rec	RPD		Limits	RPD
		Result	Result			Qual	RPD	Qual	Low	High
4,4'-DDT	18.56	0	14.2	ug/kg	77				51	146
Methoxychlor	18.56	0	13.3	ug/kg	72				54	136
Endrin ketone	18.56	0	16.3	ug/kg	88				60	129
Endrin aldehyde	18.56	0	16.6	ug/kg	89				59	132
alpha-Chlordane	18.56	0.293	17.9	ug/kg	96				39	166
gamma-Chlordane	18.56	0.2336	18.3	ug/kg	99				44	175

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q2458</u>	Analytical Method:	<u>8081B</u>
Client:	<u>CDM Smith</u>	DataFile :	<u>PD089292.D</u>

	Parameter	Spike	Sample		Units	Rec	Rec	RPD		Limits		
			Qual	RPD			Qual	Low	High	RPD		
Lab Sample ID:	Q2458-04MSD	Client Sample ID:		TP-67MSD								
	(Column 1)											
	alpha-BHC	18.54	0	18.2	ug/kg	98		3		60	144	20
	beta-BHC	18.54	0	17.4	ug/kg	94		1		54	143	20
	delta-BHC	18.54	0	19.0	ug/kg	102		6		29	151	20
	gamma-BHC (Lindane)	18.54	0	18.2	ug/kg	98		4		61	140	20
	Heptachlor	18.54	0	16.9	ug/kg	91		2		63	135	20
	Aldrin	18.54	0	18.2	ug/kg	98		1		49	139	20
	Heptachlor epoxide	18.54	0	17.7	ug/kg	95		0		41	156	20
	Endosulfan I	18.54	0	17.9	ug/kg	97		1		56	142	20
	Dieldrin	18.54	0	17.8	ug/kg	96		1		47	161	20
	4,4'-DDE	18.54	0.2077	18.0	ug/kg	95		0		55	136	20
	Endrin	18.54	0	16.9	ug/kg	91		2		57	139	20
	Endosulfan II	18.54	0	17.3	ug/kg	93		1		40	163	20
	4,4'-DDD	18.54	0	19.0	ug/kg	102		3		47	163	20
	Endosulfan sulfate	18.54	0	17.1	ug/kg	92		4		62	139	20
	4,4'-DDT	18.54	0	14.7	ug/kg	79		5		51	146	20
	Methoxychlor	18.54	0	13.8	ug/kg	74		3		54	136	20
	Endrin ketone	18.54	0	17.7	ug/kg	95		2		60	129	20
	Endrin aldehyde	18.54	0	17.3	ug/kg	93		2		59	132	20
	alpha-Chlordane	18.54	0.4784	18.2	ug/kg	96		2		39	166	20
	gamma-Chlordane	18.54	0.2856	18.2	ug/kg	98		1		44	175	20
Lab Sample ID:	Q2458-04MSD	Client Sample ID:		TP-67MSD								
	(Column 2)											
	alpha-BHC	18.54	0	18.1	ug/kg	98		3		60	144	20
	beta-BHC	18.54	0	17.9	ug/kg	97		2		54	143	20
	delta-BHC	18.54	0	17.8	ug/kg	96		5		29	151	20
	gamma-BHC (Lindane)	18.54	0	18.0	ug/kg	97		3		61	140	20
	Heptachlor	18.54	0	16.8	ug/kg	91		3		63	135	20
	Aldrin	18.54	0	18.1	ug/kg	98		2		49	139	20
	Heptachlor epoxide	18.54	0	18.3	ug/kg	99		1		41	156	20
	Endosulfan I	18.54	0	17.9	ug/kg	97		2		56	142	20
	Dieldrin	18.54	0	17.9	ug/kg	97		2		47	161	20
	4,4'-DDE	18.54	0.3449	17.9	ug/kg	97		1		55	136	20
	Endrin	18.54	0	17.2	ug/kg	93		2		57	139	20
	Endosulfan II	18.54	0	17.4	ug/kg	94		1		40	163	20
	4,4'-DDD	18.54	0	17.4	ug/kg	94		2		47	163	20
	Endosulfan sulfate	18.54	0	16.9	ug/kg	91		4		62	139	20

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q2458</u>	Analytical Method:	<u>8081B</u>
Client:	<u>CDM Smith</u>	DataFile :	<u>PD089292.D</u>

Parameter	Spike	Sample		Units	Rec	Rec		RPD		Limits	
		Result	Result			Qual	RPD	Qual	Low	High	RPD
4,4'-DDT	18.54	0	15.0	ug/kg	81		5		51	146	20
Methoxychlor	18.54	0	14.2	ug/kg	77		7		54	136	20
Endrin ketone	18.54	0	17.0	ug/kg	92		4		60	129	20
Endrin aldehyde	18.54	0	17.2	ug/kg	93		4		59	132	20
alpha-Chlordane	18.54	0.293	18.0	ug/kg	97		1		39	166	20
gamma-Chlordane	18.54	0.2336	18.4	ug/kg	99		0		44	175	20

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	Q2458	Analytical Method:	8081B
Client:	CDM Smith	Datafile :	PD089275.D

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD		Limits	
								Qual	Low	High	RPD
PB168672BS (Column 1)	alpha-BHC	16.65	18.7	ug/kg	112				84	123	
	beta-BHC	16.65	17.7	ug/kg	106				82	123	
	delta-BHC	16.65	19.2	ug/kg	115				83	126	
	gamma-BHC (Lindane)	16.65	18.6	ug/kg	112				83	125	
	Heptachlor	16.65	18.2	ug/kg	109				83	122	
	Aldrin	16.65	18.7	ug/kg	112				82	124	
	Heptachlor epoxide	16.65	18.0	ug/kg	108				83	120	
	Endosulfan I	16.65	18.3	ug/kg	110				81	124	
	Dieldrin	16.65	18.4	ug/kg	111				85	121	
	4,4'-DDE	16.65	18.4	ug/kg	111				81	123	
	Endrin	16.65	16.5	ug/kg	99				76	130	
	Endosulfan II	16.65	17.9	ug/kg	108				80	125	
	4,4'-DDD	16.65	18.7	ug/kg	112				80	131	
	Endosulfan sulfate	16.65	17.7	ug/kg	106				81	122	
	4,4'-DDT	16.65	16.3	ug/kg	98				70	129	
	Methoxychlor	16.65	16.3	ug/kg	98				60	119	
	Endrin ketone	16.65	18.3	ug/kg	110				77	132	
	Endrin aldehyde	16.65	17.8	ug/kg	107				79	124	
	alpha-Chlordane	16.65	18.1	ug/kg	109				84	120	
	gamma-Chlordane	16.65	18.5	ug/kg	111				83	122	
PB168672BS (Column 2)	alpha-BHC	16.65	18.6	ug/kg	112				84	123	
	beta-BHC	16.65	18.0	ug/kg	108				82	123	
	delta-BHC	16.65	18.3	ug/kg	110				83	126	
	gamma-BHC (Lindane)	16.65	18.4	ug/kg	111				83	125	
	Heptachlor	16.65	17.6	ug/kg	106				83	122	
	Aldrin	16.65	18.4	ug/kg	111				82	124	
	Heptachlor epoxide	16.65	19.3	ug/kg	116				83	120	
	Endosulfan I	16.65	18.4	ug/kg	111				81	124	
	Dieldrin	16.65	18.0	ug/kg	108				85	121	
	4,4'-DDE	16.65	17.9	ug/kg	108				81	123	
	Endrin	16.65	16.5	ug/kg	99				76	130	
	Endosulfan II	16.65	17.8	ug/kg	107				80	125	
	4,4'-DDD	16.65	17.9	ug/kg	108				80	131	
	Endosulfan sulfate	16.65	17.6	ug/kg	106				81	122	
	4,4'-DDT	16.65	15.9	ug/kg	95				70	129	
	Methoxychlor	16.65	16.6	ug/kg	100				60	119	
	Endrin ketone	16.65	17.8	ug/kg	107				77	132	
	Endrin aldehyde	16.65	17.6	ug/kg	106				79	124	
	alpha-Chlordane	16.65	18.2	ug/kg	109				84	120	
	gamma-Chlordane	16.65	17.9	ug/kg	108				83	122	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.: Q2458

Analytical Method: 8081B

Client: CDM Smith

Datafile : PD089330.D

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD		Limits	
								Qual	Low	High	RPD
PB168718BS (Column 1)	alpha-BHC	0.5	0.56	ug/L	111				85	130	
	beta-BHC	0.5	0.53	ug/L	106				83	126	
	delta-BHC	0.5	0.59	ug/L	117				69	141	
	gamma-BHC (Lindane)	0.5	0.55	ug/L	110				82	129	
	Heptachlor	0.5	0.54	ug/L	107				79	127	
	Aldrin	0.5	0.56	ug/L	113				79	126	
	Heptachlor epoxide	0.5	0.55	ug/L	110				81	124	
	Endosulfan I	0.5	0.56	ug/L	112				85	122	
	Dieldrin	0.5	0.57	ug/L	113				83	125	
	4,4'-DDE	0.5	0.56	ug/L	112				80	127	
	Endrin	0.5	0.49	ug/L	98				81	128	
	Endosulfan II	0.5	0.57	ug/L	113				82	123	
	4,4'-DDD	0.5	0.59	ug/L	118				77	131	
	Endosulfan sulfate	0.5	0.55	ug/L	110				76	129	
	4,4'-DDT	0.5	0.48	ug/L	96				80	133	
	Methoxychlor	0.5	0.42	ug/L	83				78	108	
	Endrin ketone	0.5	0.56	ug/L	111				80	131	
	Endrin aldehyde	0.5	0.56	ug/L	112				82	127	
	alpha-Chlordane	0.5	0.56	ug/L	111				82	125	
	gamma-Chlordane	0.5	0.57	ug/L	114				82	125	
PB168718BS (Column 2)	alpha-BHC	0.5	0.56	ug/L	112				85	130	
	beta-BHC	0.5	0.54	ug/L	109				83	126	
	delta-BHC	0.5	0.55	ug/L	110				69	141	
	gamma-BHC (Lindane)	0.5	0.56	ug/L	111				82	129	
	Heptachlor	0.5	0.52	ug/L	105				79	127	
	Aldrin	0.5	0.56	ug/L	111				79	126	
	Heptachlor epoxide	0.5	0.55	ug/L	111				81	124	
	Endosulfan I	0.5	0.55	ug/L	111				85	122	
	Dieldrin	0.5	0.55	ug/L	110				83	125	
	4,4'-DDE	0.5	0.55	ug/L	110				80	127	
	Endrin	0.5	0.48	ug/L	95				81	128	
	Endosulfan II	0.5	0.55	ug/L	110				82	123	
	4,4'-DDD	0.5	0.56	ug/L	112				77	131	
	Endosulfan sulfate	0.5	0.55	ug/L	109				76	129	
	4,4'-DDT	0.5	0.47	ug/L	95				80	133	
	Methoxychlor	0.5	0.44	ug/L	88				78	108	
	Endrin ketone	0.5	0.55	ug/L	110				80	131	
	Endrin aldehyde	0.5	0.55	ug/L	110				82	127	
	alpha-Chlordane	0.5	0.55	ug/L	109				82	125	
	gamma-Chlordane	0.5	0.55	ug/L	109				82	125	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	Q2458	Analytical Method:	8081B
Client:	CDM Smith	Datafile :	PD089331.D

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD		Limits	
								Qual	Low	High	RPD
PB168718BSD (Column 1)	alpha-BHC	0.5	0.55	ug/L	110	1			85	130	20
	beta-BHC	0.5	0.52	ug/L	104	2			83	126	20
	delta-BHC	0.5	0.57	ug/L	114	3			69	141	20
	gamma-BHC (Lindane)	0.5	0.55	ug/L	109	1			82	129	20
	Heptachlor	0.5	0.53	ug/L	106	1			79	127	20
	Aldrin	0.5	0.56	ug/L	111	2			79	126	20
	Heptachlor epoxide	0.5	0.54	ug/L	109	1			81	124	20
	Endosulfan I	0.5	0.55	ug/L	110	2			85	122	20
	Dieldrin	0.5	0.56	ug/L	112	1			83	125	20
	4,4'-DDE	0.5	0.55	ug/L	110	2			80	127	20
	Endrin	0.5	0.48	ug/L	97	1			81	128	20
	Endosulfan II	0.5	0.56	ug/L	112	1			82	123	20
	4,4'-DDD	0.5	0.58	ug/L	117	1			77	131	20
	Endosulfan sulfate	0.5	0.55	ug/L	109	1			76	129	20
	4,4'-DDT	0.5	0.47	ug/L	95	1			80	133	20
	Methoxychlor	0.5	0.42	ug/L	83	0			78	108	20
	Endrin ketone	0.5	0.55	ug/L	110	1			80	131	20
	Endrin aldehyde	0.5	0.55	ug/L	111	1			82	127	20
	alpha-Chlordane	0.5	0.55	ug/L	110	1			82	125	20
	gamma-Chlordane	0.5	0.56	ug/L	113	1			82	125	20
	alpha-BHC	0.5	0.56	ug/L	111	1			85	130	20
PB168718BSD (Column 2)	beta-BHC	0.5	0.54	ug/L	108	1			83	126	20
	delta-BHC	0.5	0.55	ug/L	110	0			69	141	20
	gamma-BHC (Lindane)	0.5	0.55	ug/L	111	0			82	129	20
	Heptachlor	0.5	0.52	ug/L	104	1			79	127	20
	Aldrin	0.5	0.55	ug/L	111	0			79	126	20
	Heptachlor epoxide	0.5	0.55	ug/L	111	0			81	124	20
	Endosulfan I	0.5	0.55	ug/L	110	1			85	122	20
	Dieldrin	0.5	0.55	ug/L	110	0			83	125	20
	4,4'-DDE	0.5	0.55	ug/L	109	1			80	127	20
	Endrin	0.5	0.48	ug/L	96	1			81	128	20
	Endosulfan II	0.5	0.54	ug/L	109	1			82	123	20
	4,4'-DDD	0.5	0.56	ug/L	112	0			77	131	20
	Endosulfan sulfate	0.5	0.55	ug/L	109	0			76	129	20
	4,4'-DDT	0.5	0.47	ug/L	94	1			80	133	20
	Methoxychlor	0.5	0.44	ug/L	88	0			78	108	20
	Endrin ketone	0.5	0.55	ug/L	111	1			80	131	20
	Endrin aldehyde	0.5	0.55	ug/L	110	0			82	127	20
	alpha-Chlordane	0.5	0.54	ug/L	108	1			82	125	20
	gamma-Chlordane	0.5	0.55	ug/L	109	0			82	125	20

4C

PESTICIDE METHOD BLANK SUMMARY

Client ID

PB168672BL

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168672BL

Lab File ID: PD089270.D

Matrix: (soil/water) Solid

Extraction: (Type) SOXH

Sulfur Cleanup: (Y/N) N

Date Extracted: 07/01/2025

Date Analyzed (1): 07/01/2025

Date Analyzed (2): 07/01/2025

Time Analyzed (1): 14:33

Time Analyzed (2): 14:33

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column (1): ZB-MR1 ID: 0.32 (mm)

GC Column (2): ZB-MR2 ID: 0.32 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED 1	DATE ANALYZED 2
PB168672BS	PB168672BS	PD089275.D	07/01/2025	07/01/2025
TP-76	Q2458-01	PD089284.D	07/01/2025	07/01/2025
TP-55	Q2458-02	PD089285.D	07/01/2025	07/01/2025
TP-68	Q2458-03	PD089286.D	07/01/2025	07/01/2025
TP-67	Q2458-04	PD089290.D	07/01/2025	07/01/2025
TP-67MS	Q2458-04MS	PD089291.D	07/01/2025	07/01/2025
TP-67MSD	Q2458-04MSD	PD089292.D	07/01/2025	07/01/2025
TP-66	Q2458-05	PD089293.D	07/01/2025	07/01/2025
TP-60	Q2458-06	PD089294.D	07/01/2025	07/01/2025
TP-62	Q2458-07	PD089295.D	07/01/2025	07/01/2025
TP-63	Q2458-08	PD089296.D	07/01/2025	07/01/2025
TP-59	Q2458-09	PD089297.D	07/01/2025	07/01/2025

COMMENTS: _____

4C

PESTICIDE METHOD BLANK SUMMARY

Client ID

PB168718BL

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168718BL

Lab File ID: PD089329.D

Matrix: (soil/water) WATER

Extraction: (Type) SEPF

Sulfur Cleanup: (Y/N) N

Date Extracted: 07/03/2025

Date Analyzed (1): 07/03/2025

Date Analyzed (2): 07/03/2025

Time Analyzed (1): 13:55

Time Analyzed (2): 13:55

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column (1): ZB-MR1 ID: 0.32 (mm)

GC Column (2): ZB-MR2 ID: 0.32 (mm)

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

EPA SAMPLE NO.	LAB SAMPLE ID	LAB FILE ID	DATE ANALYZED 1	DATE ANALYZED 2
PB168718BS	PB168718BS	PD089330.D	07/03/2025	07/03/2025
PB168718BSD	PB168718BSD	PD089331.D	07/03/2025	07/03/2025
FB-06272025	Q2458-10	PD089332.D	07/03/2025	07/03/2025

COMMENTS: _____



SAMPLE DATA

Report of Analysis

Client:	CDM Smith	Date Collected:	06/26/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-76	SDG No.:	Q2458
Lab Sample ID:	Q2458-01	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	90.6 Decanted:
Sample Wt/Vol:	30.04 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089284.D	1	07/01/25 08:30	07/01/25 18:26	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.90	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.90	ug/kg
319-86-8	delta-BHC	0.43	U	0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	0.15	U	0.15	1.90	ug/kg
76-44-8	Heptachlor	0.13	U	0.13	1.90	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.90	ug/kg
959-98-8	Endosulfan I	0.15	U	0.15	1.90	ug/kg
60-57-1	Dieldrin	1.90	P	0.15	1.90	ug/kg
72-55-9	4,4-DDE	0.15	U	0.15	1.90	ug/kg
72-20-8	Endrin	0.66	J	0.15	1.90	ug/kg
33213-65-9	Endosulfan II	0.32	U	0.32	1.90	ug/kg
72-54-8	4,4-DDD	0.17	U	0.17	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.90	ug/kg
50-29-3	4,4-DDT	1.20	JP	0.15	1.90	ug/kg
72-43-5	Methoxychlor	0.41	U	0.41	1.90	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	0.41	U	0.41	1.90	ug/kg
5103-71-9	alpha-Chlordane	0.13	U	0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	0.17	U	0.17	1.90	ug/kg
8001-35-2	Toxaphene	6.00	U	6.00	36.4	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	13.9		20 - 144	70%	SPK: 20
877-09-8	Tetrachloro-m-xylene	17.8		19 - 148	89%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/26/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-76	SDG No.:	Q2458
Lab Sample ID:	Q2458-01	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	90.6 Decanted:
Sample Wt/Vol:	30.04 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089284.D	1	07/01/25 08:30	07/01/25 18:26	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089284.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 18:26
 Operator : AR\AJ
 Sample : Q2458-01
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-76

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/02/2025
 Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:49:42 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.879	47188302	301.5E6	16.539	17.787m
28)	SA Decachlor...	9.072	8.070	54771579	259.7E6	13.945m	13.134
Target Compounds							
13)	MA Dieldrin	6.351	5.511	6458897	116.9E6	1.346m	5.038m#
14)	MA Endrin	6.576	5.778	7395274	35884222	1.793	1.653m
17)	MA 4,4'-DDT	7.025	6.178	5295917	66983031	1.398	3.303 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089284.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 18:26
Operator : AR\AJ
Sample : Q2458-01
Misc :
ALS Vial : 19 Sample Multiplier: 1

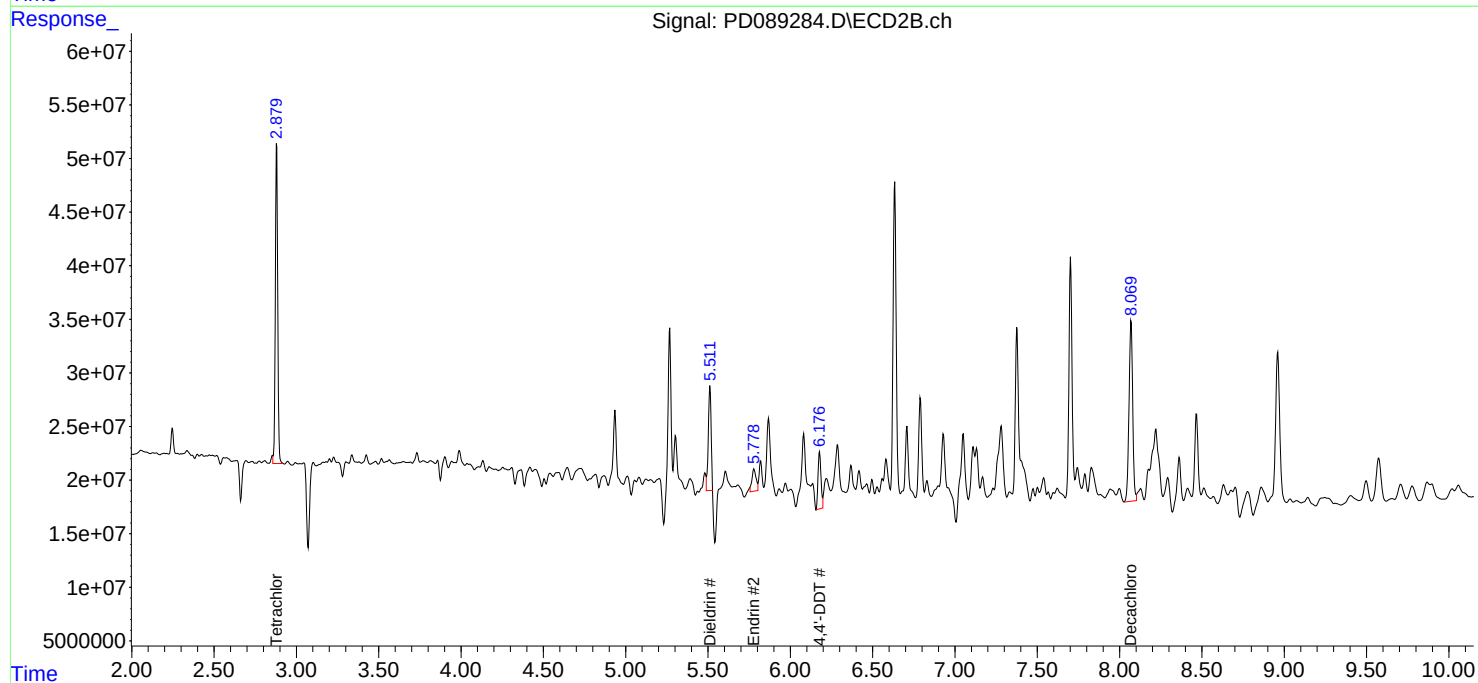
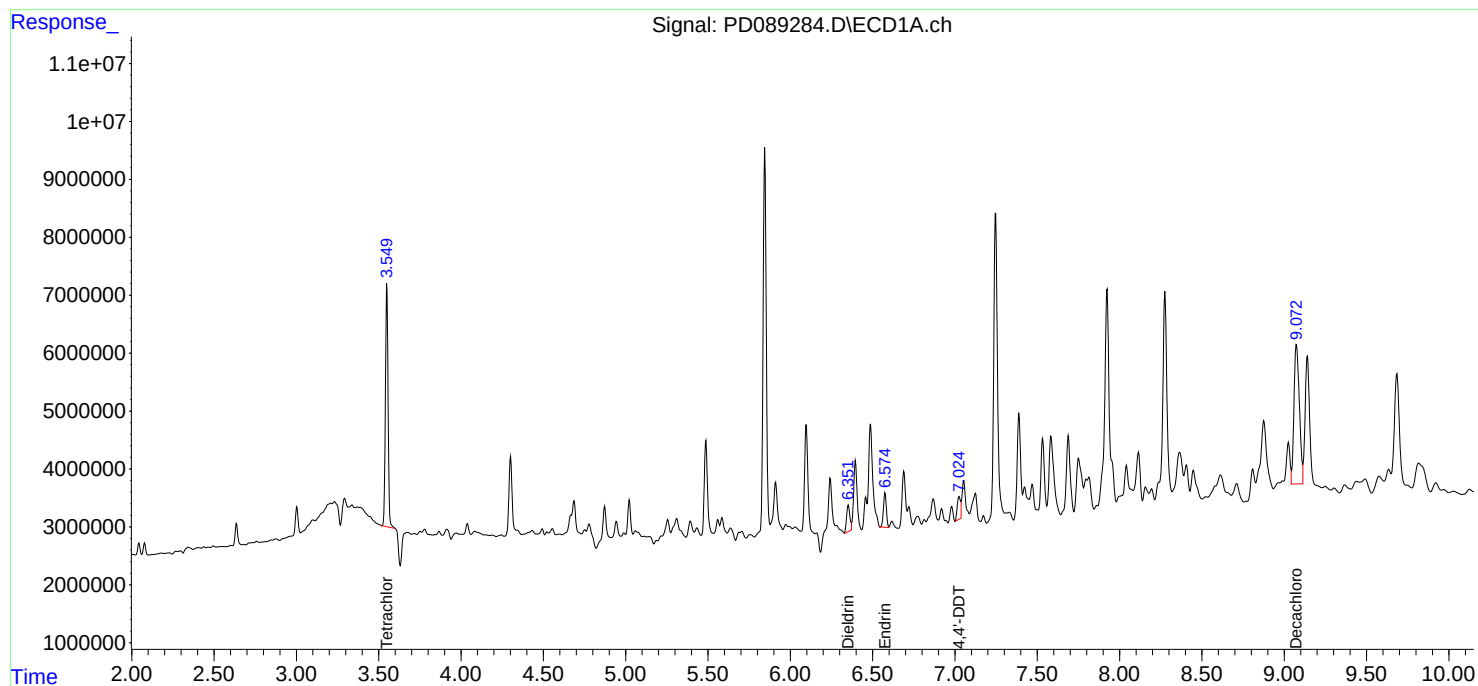
Instrument :
ECD_D
ClientSampleId :
TP-76

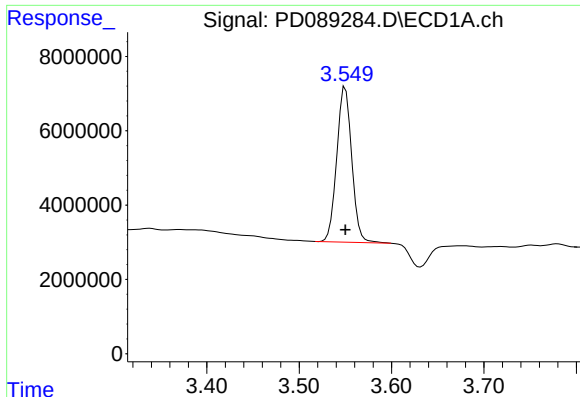
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:49:42 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





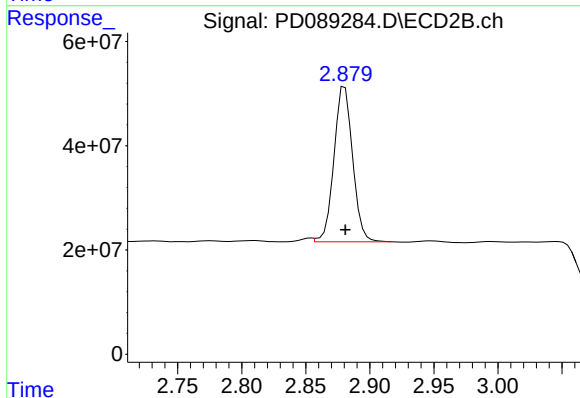
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
 Delta R.T.: 0.000 min
 Response: 47188302
 Conc: 16.54 ng/ml

Instrument :
 ECD_D
 ClientSampleId :
 TP-76

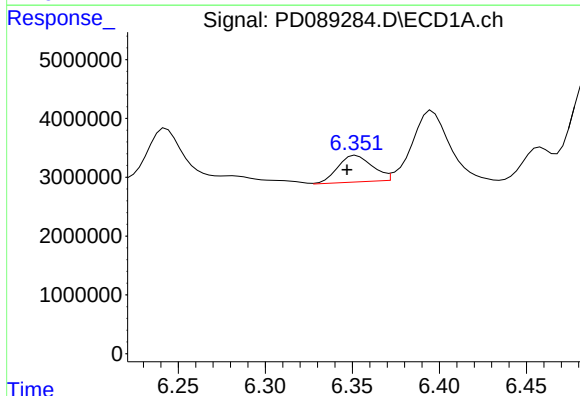
Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025



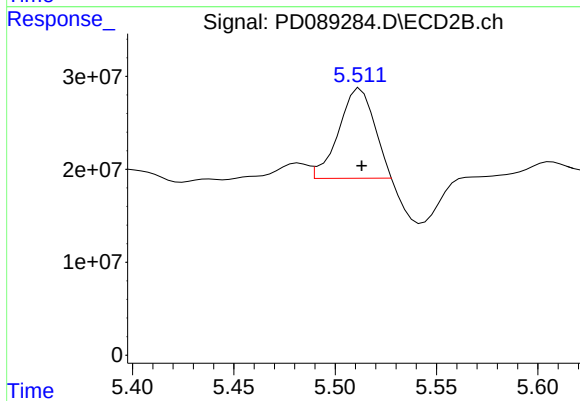
#1 Tetrachloro-m-xylene

R.T.: 2.879 min
 Delta R.T.: -0.002 min
 Response: 301544783
 Conc: 17.79 ng/ml m



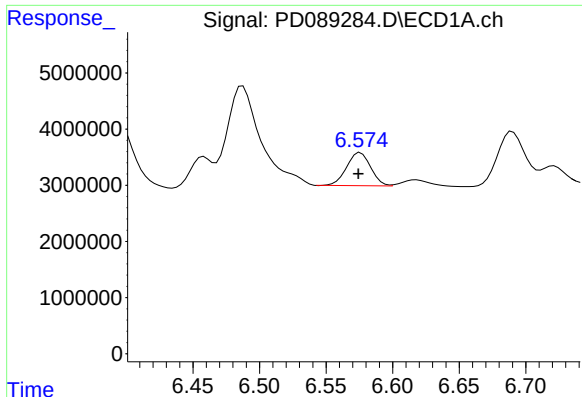
#13 Dieldrin

R.T.: 6.351 min
 Delta R.T.: 0.004 min
 Response: 6458897
 Conc: 1.35 ng/ml m



#13 Dieldrin

R.T.: 5.511 min
 Delta R.T.: -0.002 min
 Response: 116924415
 Conc: 5.04 ng/ml m



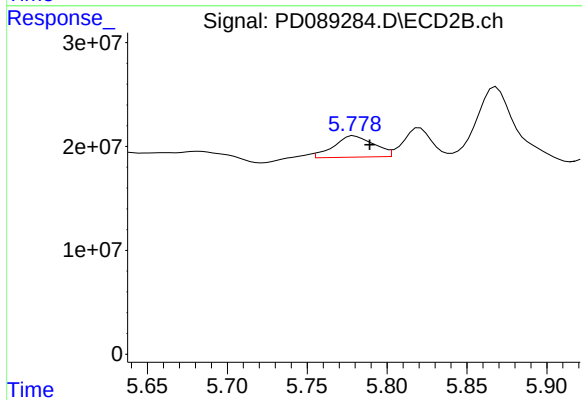
#14 Endrin

R.T.: 6.576 min
 Delta R.T.: 0.001 min
 Response: 7395274
 Conc: 1.79 ng/ml

Instrument :
 ECD_D
 ClientSampleId :
 TP-76

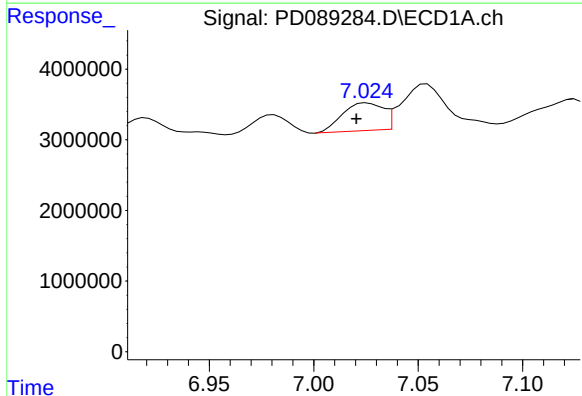
Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025



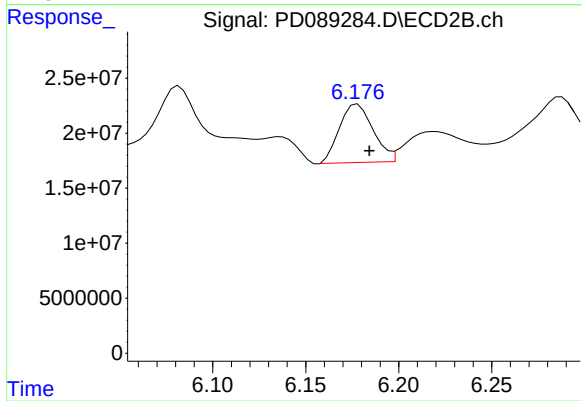
#14 Endrin

R.T.: 5.778 min
 Delta R.T.: -0.011 min
 Response: 35884222
 Conc: 1.65 ng/ml m



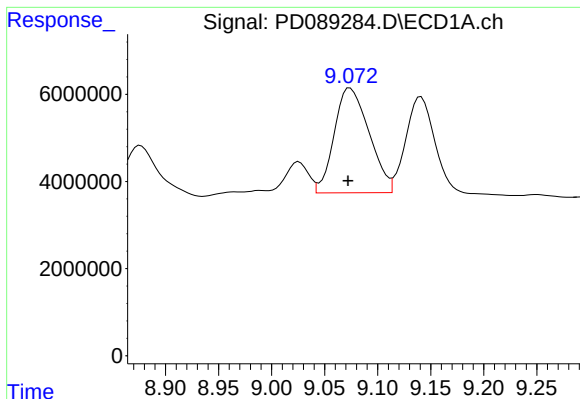
#17 4,4'-DDT

R.T.: 7.025 min
 Delta R.T.: 0.005 min
 Response: 5295917
 Conc: 1.40 ng/ml



#17 4,4'-DDT

R.T.: 6.178 min
 Delta R.T.: -0.006 min
 Response: 66983031
 Conc: 3.30 ng/ml



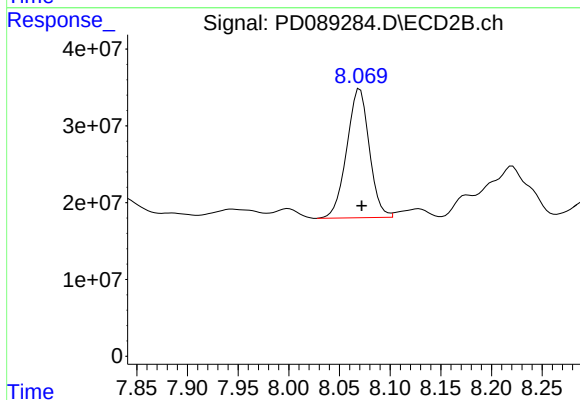
#28 Decachlorobiphenyl

R.T.: 9.072 min
Delta R.T.: 0.000 min
Response: 54771579
Conc: 13.95 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-76

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



#28 Decachlorobiphenyl

R.T.: 8.070 min
Delta R.T.: -0.002 min
Response: 259659292
Conc: 13.13 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/26/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-55	SDG No.:	Q2458
Lab Sample ID:	Q2458-02	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	91.4 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089285.D	1	07/01/25 08:30	07/01/25 18:40	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.90	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.90	ug/kg
319-86-8	delta-BHC	0.43	U	0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	0.15	U	0.15	1.90	ug/kg
76-44-8	Heptachlor	0.13	U	0.13	1.90	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.90	ug/kg
959-98-8	Endosulfan I	0.15	U	0.15	1.90	ug/kg
60-57-1	Dieldrin	0.15	U	0.15	1.90	ug/kg
72-55-9	4,4-DDE	0.15	U	0.15	1.90	ug/kg
72-20-8	Endrin	0.15	U	0.15	1.90	ug/kg
33213-65-9	Endosulfan II	0.32	U	0.32	1.90	ug/kg
72-54-8	4,4-DDD	0.16	U	0.16	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.90	ug/kg
50-29-3	4,4-DDT	0.15	U	0.15	1.90	ug/kg
72-43-5	Methoxychlor	0.40	U	0.40	1.90	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	0.40	U	0.40	1.90	ug/kg
5103-71-9	alpha-Chlordane	0.13	U	0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	0.16	U	0.16	1.90	ug/kg
8001-35-2	Toxaphene	5.90	U	5.90	36.1	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	9.41		20 - 144	47%	SPK: 20
877-09-8	Tetrachloro-m-xylene	11.8		19 - 148	59%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/26/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-55	SDG No.:	Q2458
Lab Sample ID:	Q2458-02	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	91.4 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089285.D	1	07/01/25 08:30	07/01/25 18:40	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089285.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 18:40
Operator : AR\AJ
Sample : Q2458-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
TP-55

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:49:57 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	31512300	200.2E6	11.045	11.806
28)	SA Decachlor...	9.073	8.072	36952271	147.3E6	9.408	7.449

Target Compounds

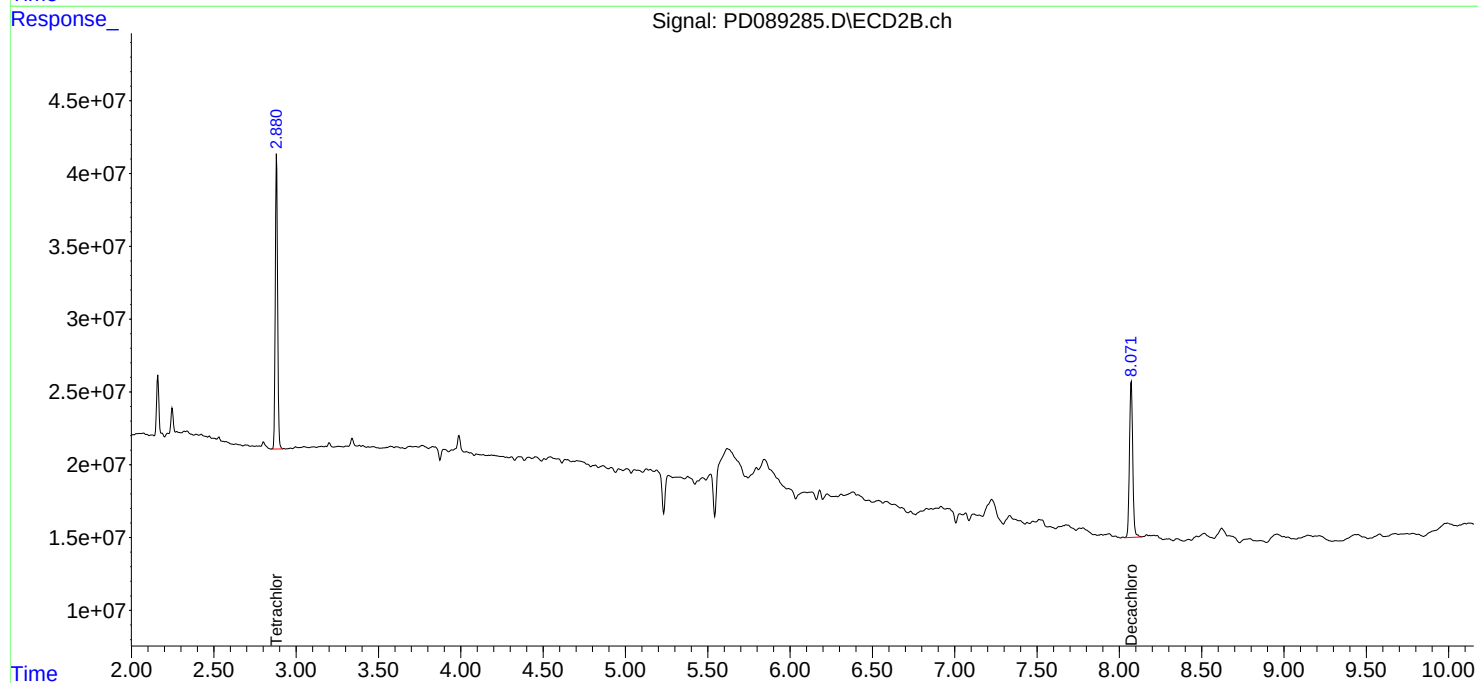
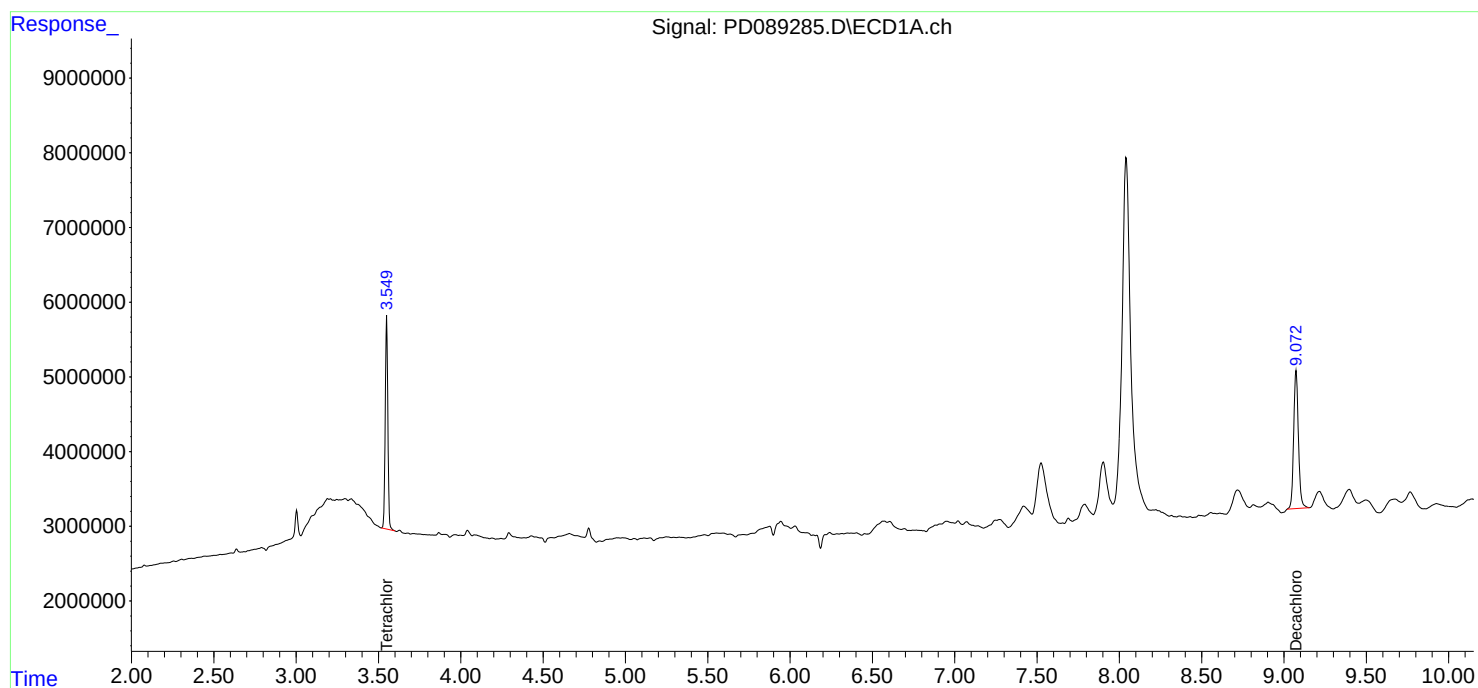
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

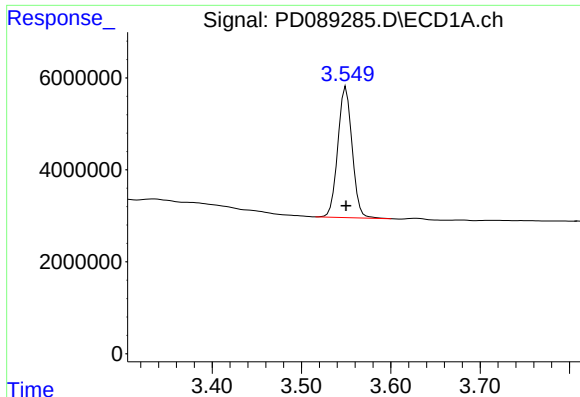
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089285.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 18:40
Operator : AR\AJ
Sample : Q2458-02
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
TP-55

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:49:57 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

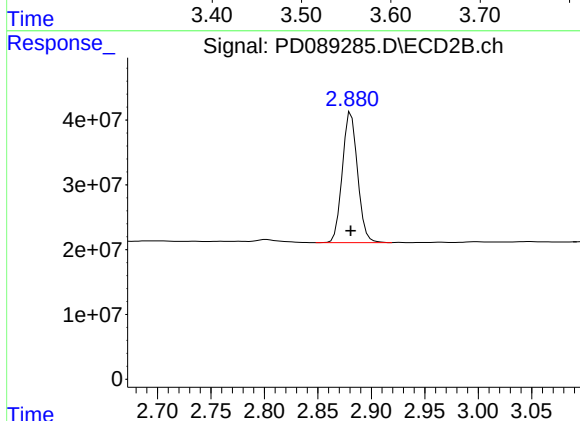




#1 Tetrachloro-m-xylene

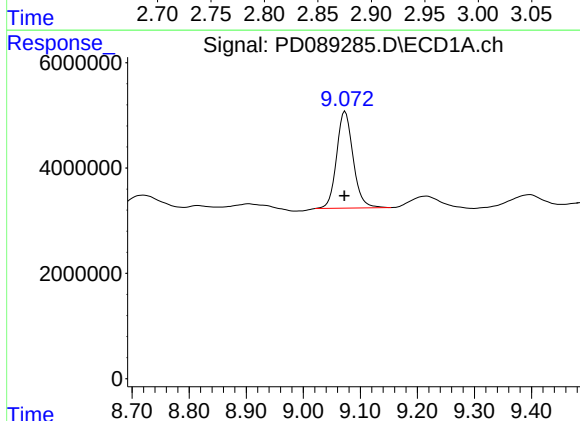
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 31512300
Conc: 11.04 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-55



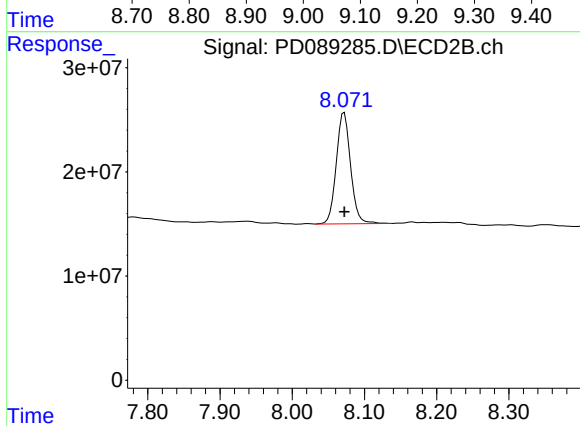
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 200155158
Conc: 11.81 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.002 min
Response: 36952271
Conc: 9.41 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 147262407
Conc: 7.45 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-68	SDG No.:	Q2458
Lab Sample ID:	Q2458-03	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	92.3 Decanted:
Sample Wt/Vol:	30.01 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089286.D	1	07/01/25 08:30	07/01/25 18:53	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.80	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.80	ug/kg
319-86-8	delta-BHC	0.42	U	0.42	1.80	ug/kg
58-89-9	gamma-BHC (Lindane)	0.15	U	0.15	1.80	ug/kg
76-44-8	Heptachlor	0.13	U	0.13	1.80	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.80	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.80	ug/kg
959-98-8	Endosulfan I	0.15	U	0.15	1.80	ug/kg
60-57-1	Dieldrin	0.15	U	0.15	1.80	ug/kg
72-55-9	4,4-DDE	0.15	U	0.15	1.80	ug/kg
72-20-8	Endrin	0.15	U	0.15	1.80	ug/kg
33213-65-9	Endosulfan II	0.31	U	0.31	1.80	ug/kg
72-54-8	4,4-DDD	0.16	U	0.16	1.80	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.80	ug/kg
50-29-3	4,4-DDT	0.15	U	0.15	1.80	ug/kg
72-43-5	Methoxychlor	0.40	U	0.40	1.80	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.80	ug/kg
7421-93-4	Endrin aldehyde	0.40	U	0.40	1.80	ug/kg
5103-71-9	alpha-Chlordane	0.13	U	0.13	1.80	ug/kg
5103-74-2	gamma-Chlordane	0.16	U	0.16	1.80	ug/kg
8001-35-2	Toxaphene	5.90	U	5.90	35.7	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	13.1		20 - 144	66%	SPK: 20
877-09-8	Tetrachloro-m-xylene	17.9		19 - 148	89%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-68	SDG No.:	Q2458
Lab Sample ID:	Q2458-03	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	92.3 Decanted:
Sample Wt/Vol:	30.01 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089286.D	1	07/01/25 08:30	07/01/25 18:53	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089286.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 18:53
 Operator : AR\AJ
 Sample : Q2458-03
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-68

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:50:10 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.550	2.879	46258267	302.6E6	16.213	17.852m
28) SA Decachlor...	9.072	8.070	51575801	216.0E6	13.132m	10.926

Target Compounds

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089286.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 18:53
Operator : AR\AJ
Sample : Q2458-03
Misc :
ALS Vial : 21 Sample Multiplier: 1

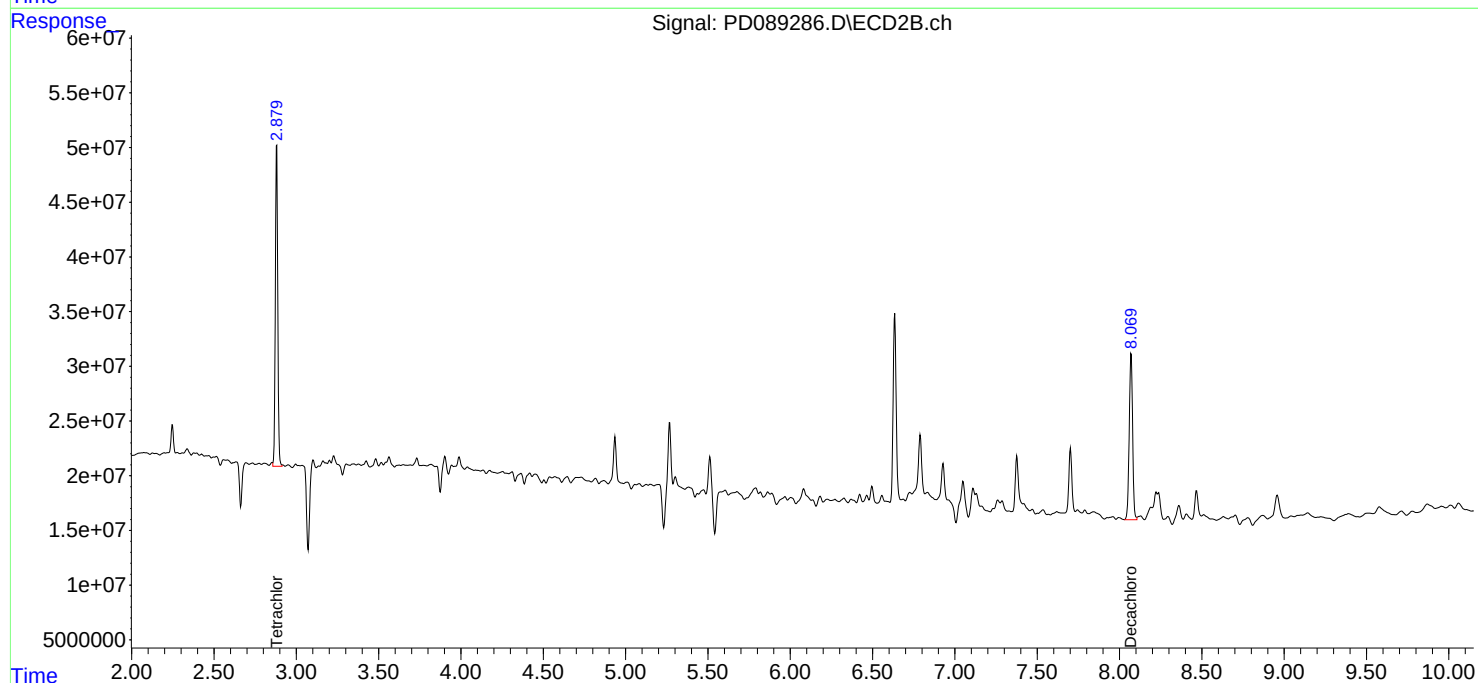
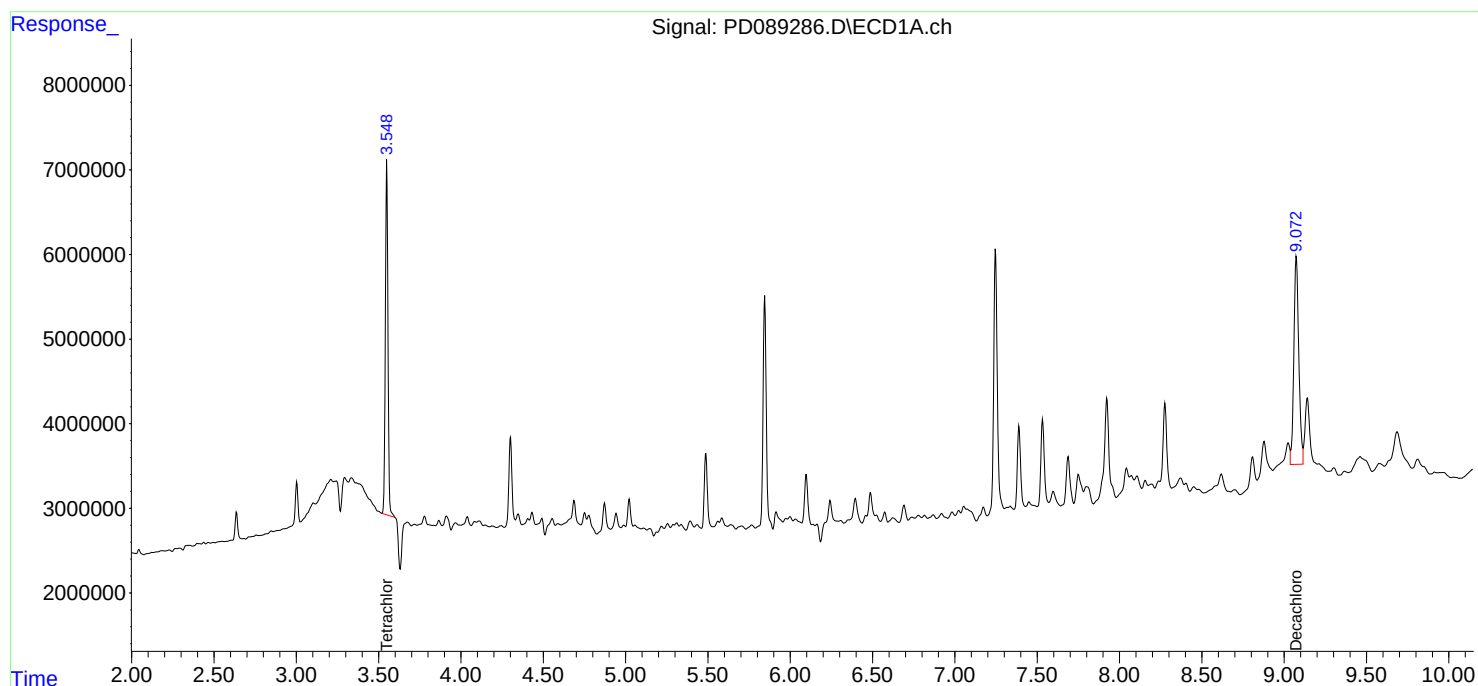
Instrument :
ECD_D
ClientSampleId :
TP-68

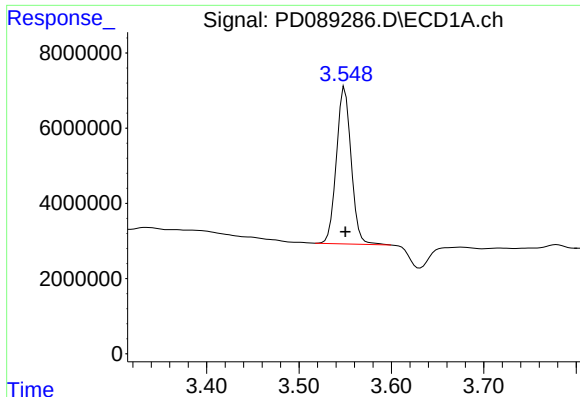
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:50:10 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





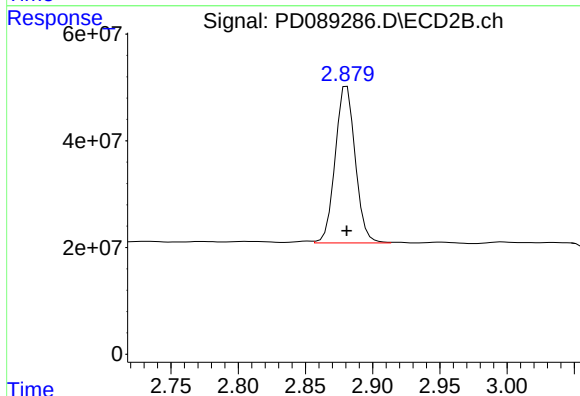
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
 Delta R.T.: 0.000 min
 Response: 46258267
 Conc: 16.21 ng/ml

Instrument :
 ECD_D
 ClientSampleId :
 TP-68

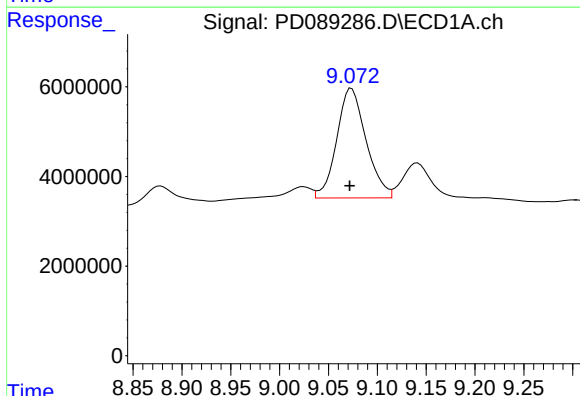
Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025



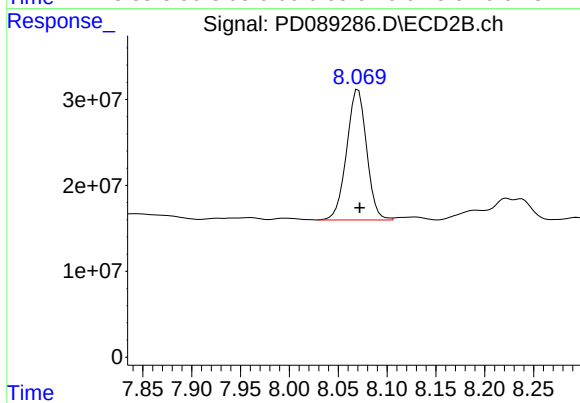
#1 Tetrachloro-m-xylene

R.T.: 2.879 min
 Delta R.T.: -0.001 min
 Response: 302645220
 Conc: 17.85 ng/ml m



#28 Decachlorobiphenyl

R.T.: 9.072 min
 Delta R.T.: 0.000 min
 Response: 51575801
 Conc: 13.13 ng/ml m



#28 Decachlorobiphenyl

R.T.: 8.070 min
 Delta R.T.: -0.002 min
 Response: 216010092
 Conc: 10.93 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67	SDG No.:	Q2458
Lab Sample ID:	Q2458-04	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7 Decanted:
Sample Wt/Vol:	30.06 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089290.D	1	07/01/25 08:30	07/01/25 20:15	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.90	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.90	ug/kg
319-86-8	delta-BHC	0.43	U	0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	0.16	U	0.16	1.90	ug/kg
76-44-8	Heptachlor	0.13	U	0.13	1.90	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.90	ug/kg
959-98-8	Endosulfan I	0.16	U	0.16	1.90	ug/kg
60-57-1	Dieldrin	0.16	U	0.16	1.90	ug/kg
72-55-9	4,4-DDE	0.34	JP	0.16	1.90	ug/kg
72-20-8	Endrin	0.16	U	0.16	1.90	ug/kg
33213-65-9	Endosulfan II	0.32	U	0.32	1.90	ug/kg
72-54-8	4,4-DDD	0.17	U	0.17	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.90	ug/kg
50-29-3	4,4-DDT	0.16	U	0.16	1.90	ug/kg
72-43-5	Methoxychlor	0.41	U	0.41	1.90	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	0.41	U	0.41	1.90	ug/kg
5103-71-9	alpha-Chlordane	0.48	JP	0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	0.29	J	0.17	1.90	ug/kg
8001-35-2	Toxaphene	6.00	U	6.00	36.7	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	15.0		20 - 144	75%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.0		19 - 148	100%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67	SDG No.:	Q2458
Lab Sample ID:	Q2458-04	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7 Decanted:
Sample Wt/Vol:	30.06 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089290.D	1	07/01/25 08:30	07/01/25 20:15	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089290.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 20:15
 Operator : AR\AJ
 Sample : Q2458-04
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-67

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/02/2025
 Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:51:23 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.551	2.880	54454459	339.4E6	19.086	20.019m
28) SA Decachlor...	9.074	8.072	58649677	295.7E6	14.933	14.956
Target Compounds						
10) B gamma-Chl...	5.944	5.125	3682776	15008798	0.769m	0.627m
11) B alpha-Chl...	6.031	5.189	6243612	18011582	1.290	0.787m#
12) B 4,4'-DDE	6.196	5.374	2459007	21389348	0.564	0.933m#
17) MA 4,4'-DDT	7.020	6.183	1329797	21703639	0.351m	1.070 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089290.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 20:15
Operator : AR\AJ
Sample : Q2458-04
Misc :
ALS Vial : 22 Sample Multiplier: 1

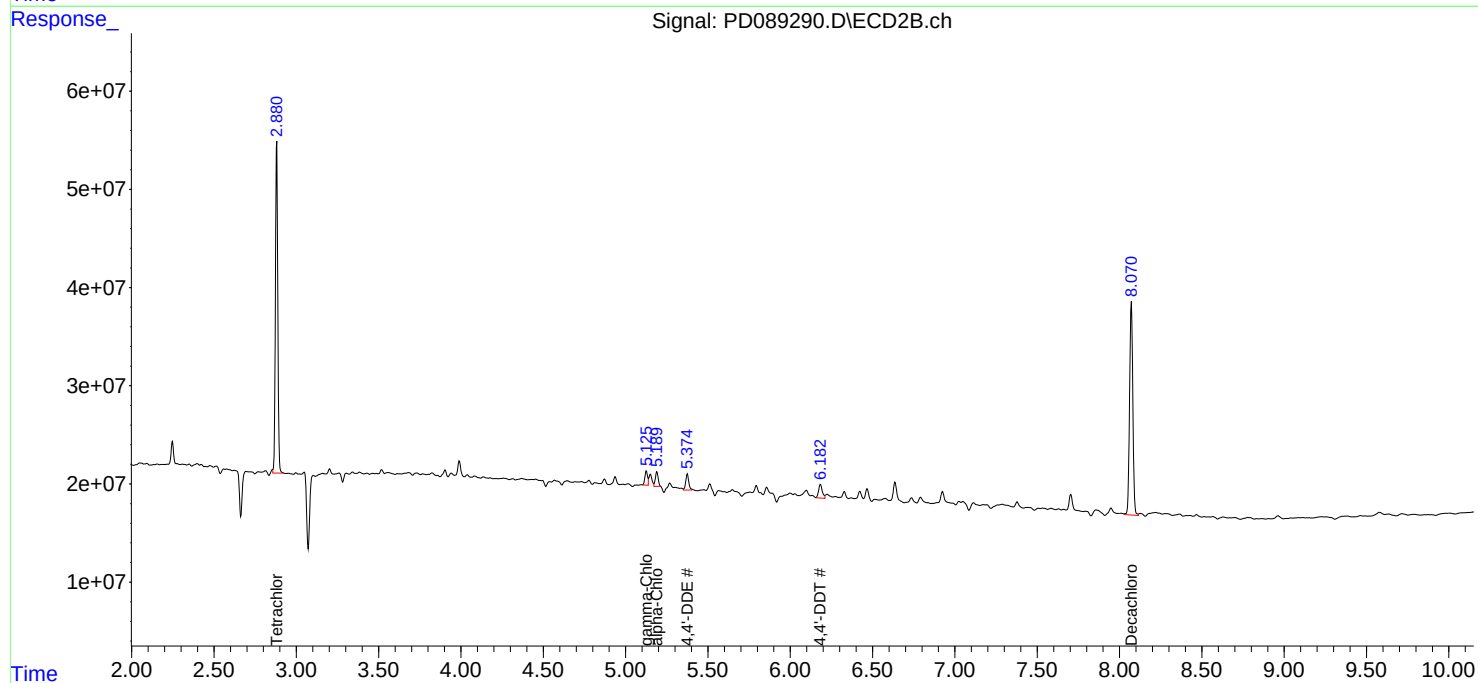
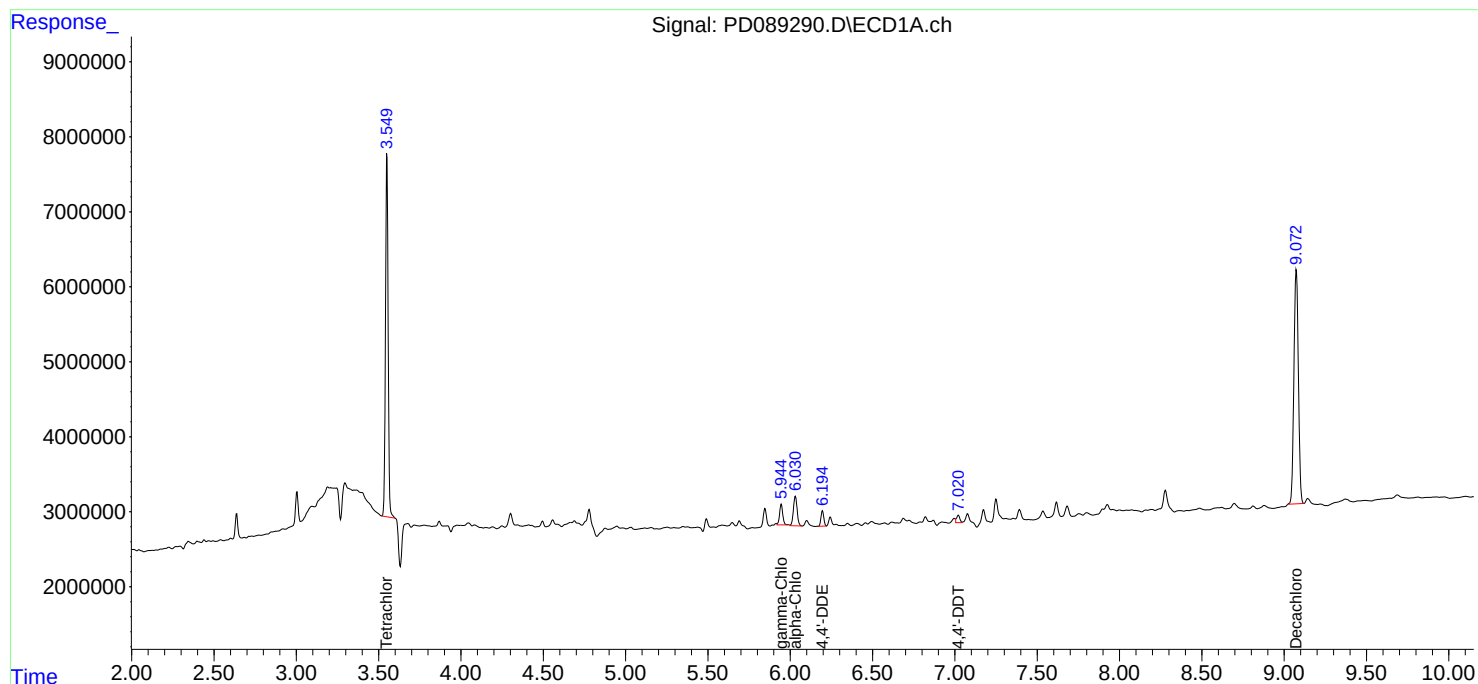
Instrument :
ECD_D
ClientSampleId :
TP-67

Manual Integrations
APPROVED

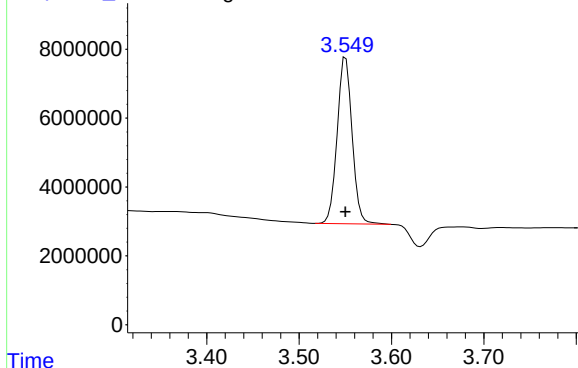
Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:51:23 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Response_ Signal: PD089290.D\ECD1A.ch



#1 Tetrachloro-m-xylene

R.T.: 3.551 min
 Delta R.T.: 0.000 min
 Response: 54454459
 Conc: 19.09 ng/ml

Instrument :

ECD_D

ClientSampleId :

TP-67

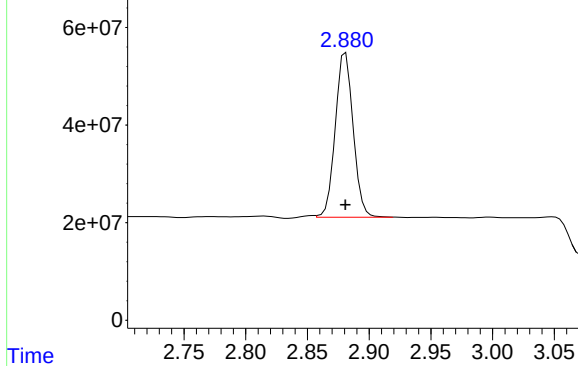
Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

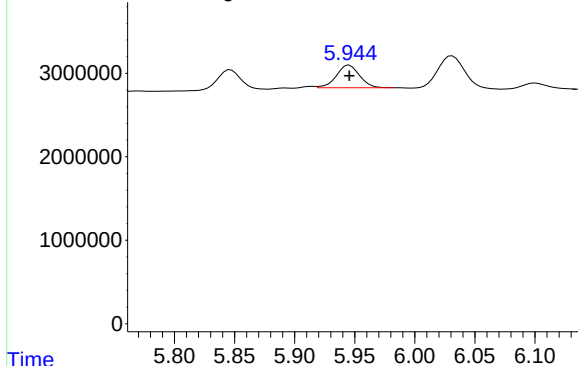
Time Response_ Signal: PD089290.D\ECD2B.ch



#1 Tetrachloro-m-xylene

R.T.: 2.880 min
 Delta R.T.: -0.001 min
 Response: 339387727
 Conc: 20.02 ng/ml m

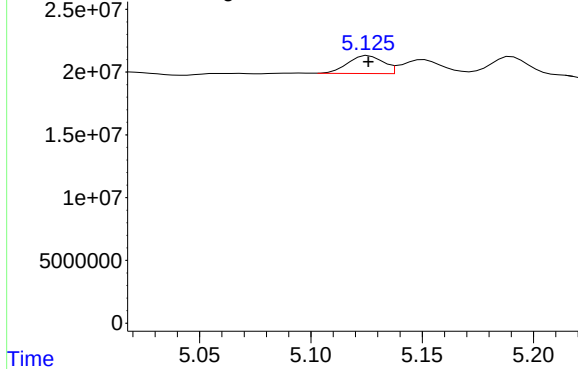
Time Response_ Signal: PD089290.D\ECD1A.ch



#10 gamma-Chlordane

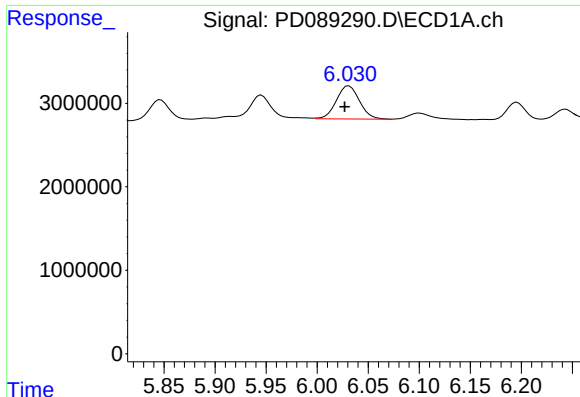
R.T.: 5.944 min
 Delta R.T.: -0.001 min
 Response: 3682776
 Conc: 0.77 ng/ml m

Time Response_ Signal: PD089290.D\ECD2B.ch



#10 gamma-Chlordane

R.T.: 5.125 min
 Delta R.T.: -0.001 min
 Response: 15008798
 Conc: 0.63 ng/ml m



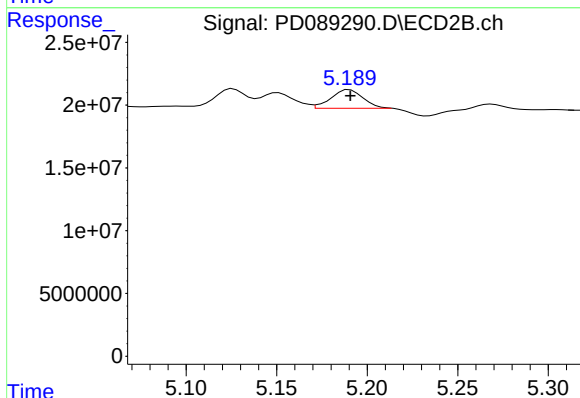
#11 alpha-Chlordane

R.T.: 6.031 min
Delta R.T.: 0.004 min
Response: 6243612
Conc: 1.29 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-67

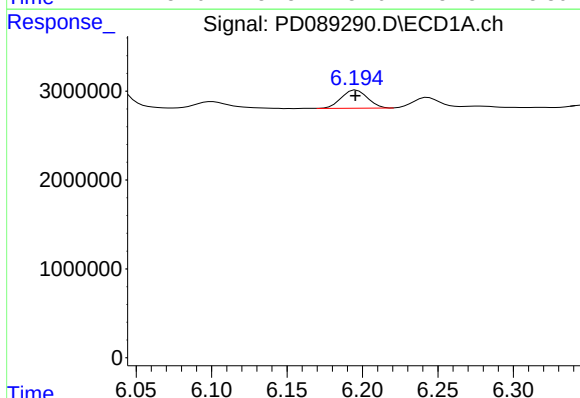
Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



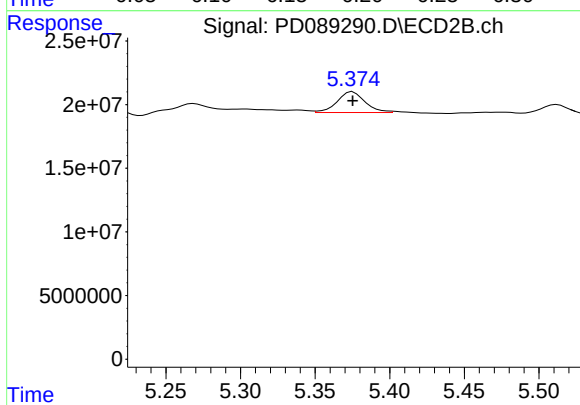
#11 alpha-Chlordane

R.T.: 5.189 min
Delta R.T.: -0.002 min
Response: 18011582
Conc: 0.79 ng/ml m



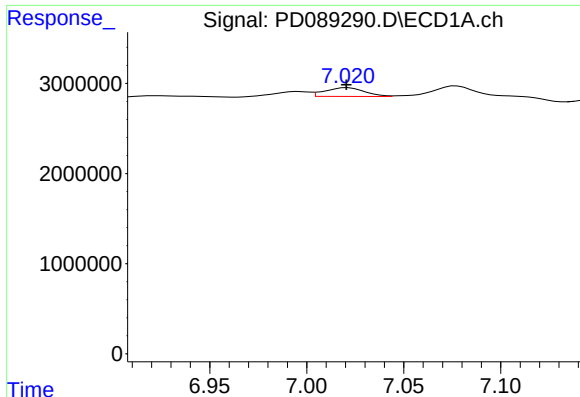
#12 4,4'-DDE

R.T.: 6.196 min
Delta R.T.: 0.000 min
Response: 2459007
Conc: 0.56 ng/ml



#12 4,4'-DDE

R.T.: 5.374 min
Delta R.T.: -0.001 min
Response: 21389348
Conc: 0.93 ng/ml m



#17 4,4'-DDT

R.T.: 7.020 min
Delta R.T.: 0.000 min
Response: 1329797
Conc: 0.35 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-67

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025

#17 4,4'-DDT

R.T.: 6.183 min
Delta R.T.: 0.000 min
Response: 21703639
Conc: 1.07 ng/ml

#28 Decachlorobiphenyl

R.T.: 9.074 min
Delta R.T.: 0.002 min
Response: 58649677
Conc: 14.93 ng/ml

#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 295669920
Conc: 14.96 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-66	SDG No.:	Q2458
Lab Sample ID:	Q2458-05	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	88.3 Decanted:
Sample Wt/Vol:	30.06 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089293.D	1	07/01/25 08:30	07/01/25 20:57	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.15	U	0.15	1.90	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.90	ug/kg
319-86-8	delta-BHC	0.44	U	0.44	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	0.16	U	0.16	1.90	ug/kg
76-44-8	Heptachlor	0.14	U	0.14	1.90	ug/kg
309-00-2	Aldrin	0.14	U	0.14	1.90	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.90	ug/kg
959-98-8	Endosulfan I	0.16	U	0.16	1.90	ug/kg
60-57-1	Dieldrin	0.16	U	0.16	1.90	ug/kg
72-55-9	4,4-DDE	0.16	U	0.16	1.90	ug/kg
72-20-8	Endrin	0.16	U	0.16	1.90	ug/kg
33213-65-9	Endosulfan II	0.33	U	0.33	1.90	ug/kg
72-54-8	4,4-DDD	0.17	U	0.17	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	0.15	U	0.15	1.90	ug/kg
50-29-3	4,4-DDT	0.40	JP	0.16	1.90	ug/kg
72-43-5	Methoxychlor	0.42	U	0.42	1.90	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	0.42	U	0.42	1.90	ug/kg
5103-71-9	alpha-Chlordane	0.14	U	0.14	1.90	ug/kg
5103-74-2	gamma-Chlordane	0.17	U	0.17	1.90	ug/kg
8001-35-2	Toxaphene	6.10	U	6.10	37.3	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	10.6		20 - 144	53%	SPK: 20
877-09-8	Tetrachloro-m-xylene	12.6		19 - 148	63%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-66	SDG No.:	Q2458
Lab Sample ID:	Q2458-05	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	88.3 Decanted:
Sample Wt/Vol:	30.06 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089293.D	1	07/01/25 08:30	07/01/25 20:57	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089293.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 20:57
Operator : AR\AJ
Sample : Q2458-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
TP-66

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:00 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.551	2.881	33039023	213.6E6	11.580	12.598
28) SA Decachlor...	9.073	8.072	41602422	206.3E6	10.592	10.434
Target Compounds						
17) MA 4,4'-DDT	7.020	6.179	2447978	21604807	0.646m	1.065m#

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089293.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 20:57
Operator : AR\AJ
Sample : Q2458-05
Misc :
ALS Vial : 25 Sample Multiplier: 1

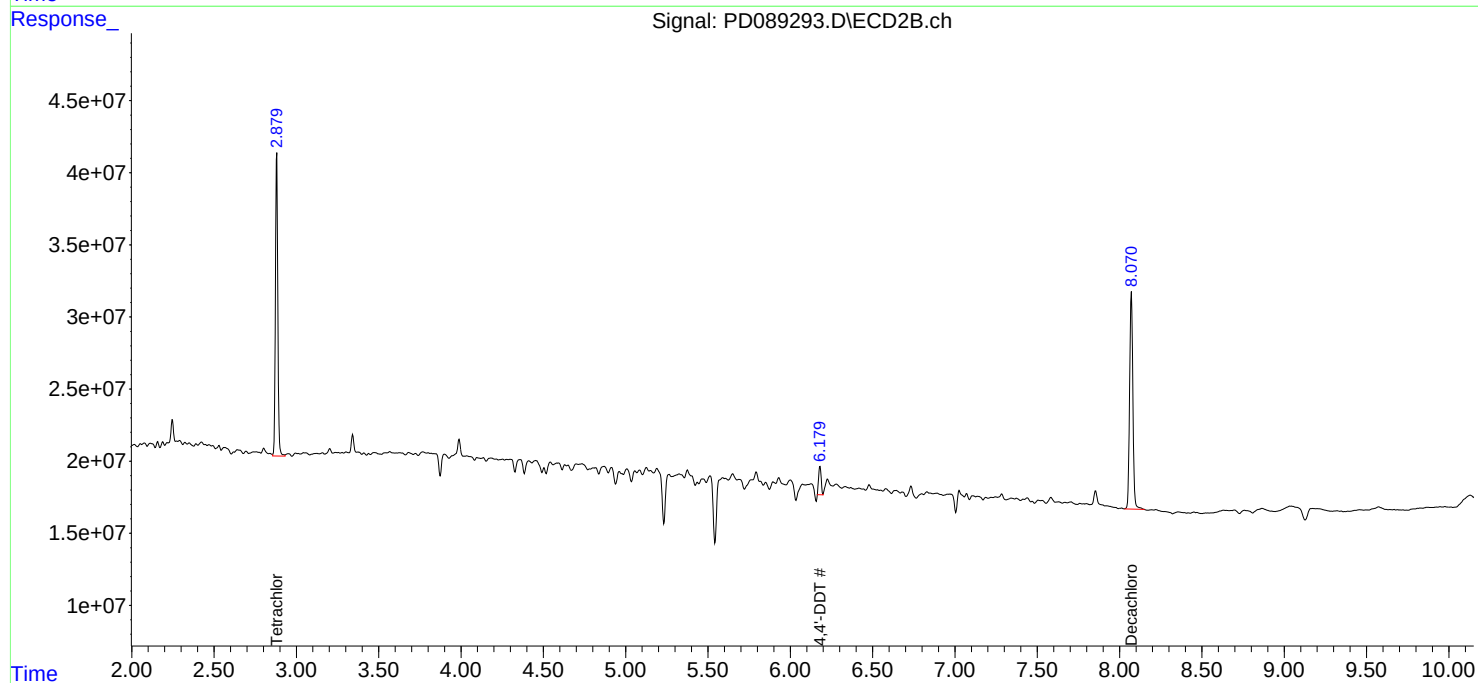
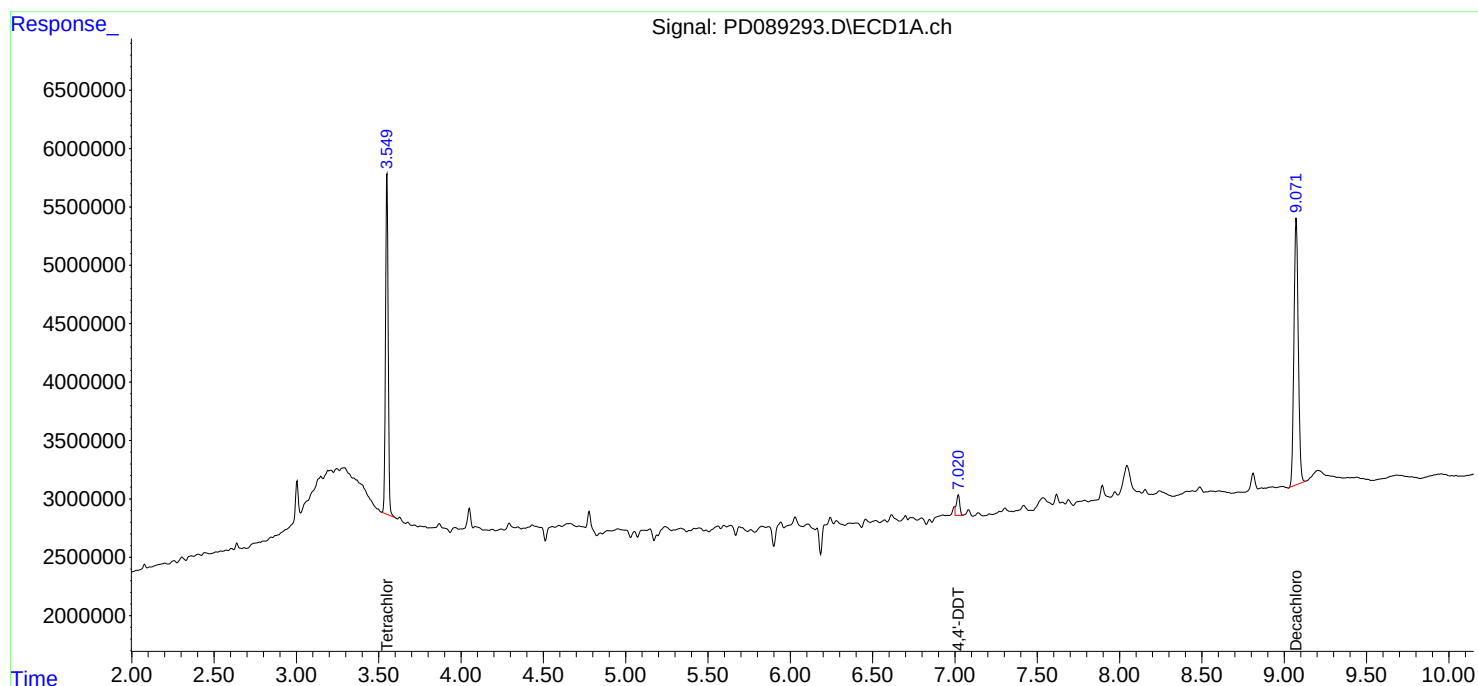
Instrument :
ECD_D
ClientSampleId :
TP-66

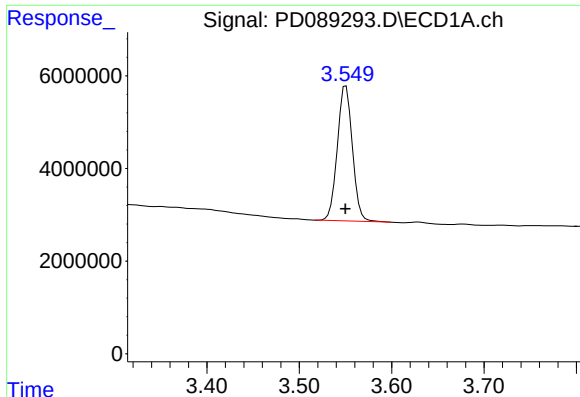
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:00 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





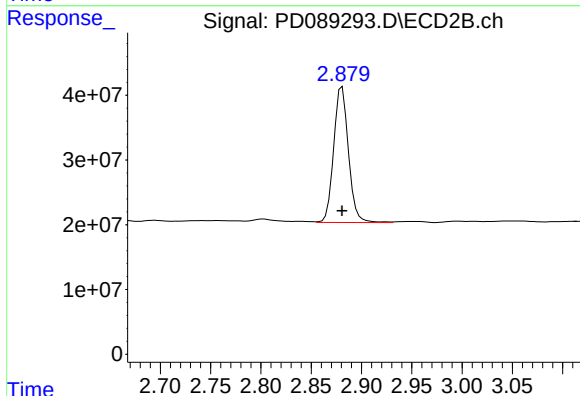
#1 Tetrachloro-m-xylene

R.T.: 3.551 min
 Delta R.T.: 0.000 min
 Response: 33039023
 Conc: 11.58 ng/ml

Instrument :
 ECD_D
 ClientSampleId :
 TP-66

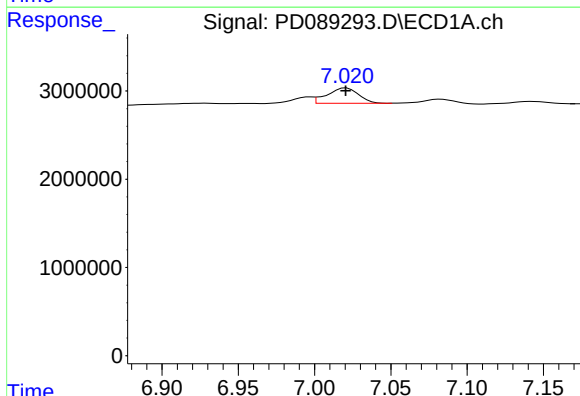
Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025



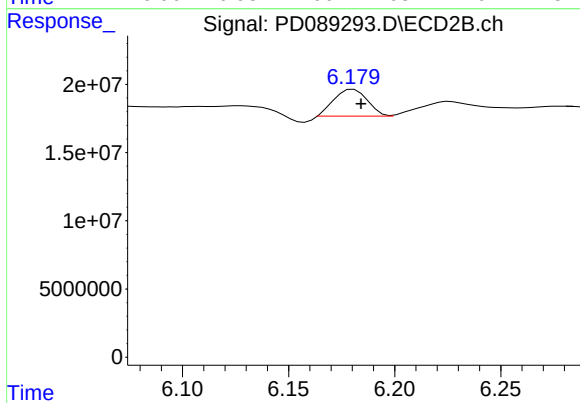
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
 Delta R.T.: 0.000 min
 Response: 213570595
 Conc: 12.60 ng/ml



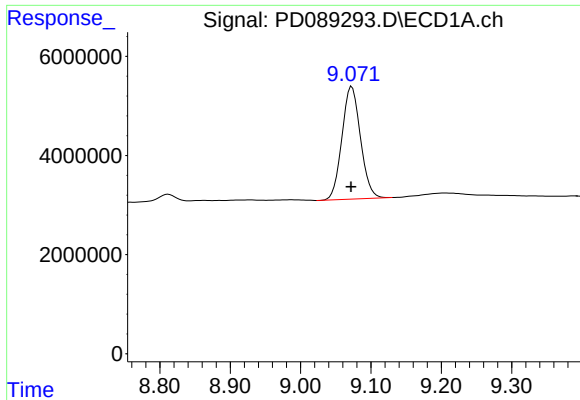
#17 4,4'-DDT

R.T.: 7.020 min
 Delta R.T.: 0.000 min
 Response: 2447978
 Conc: 0.65 ng/ml m



#17 4,4'-DDT

R.T.: 6.179 min
 Delta R.T.: -0.005 min
 Response: 21604807
 Conc: 1.07 ng/ml m



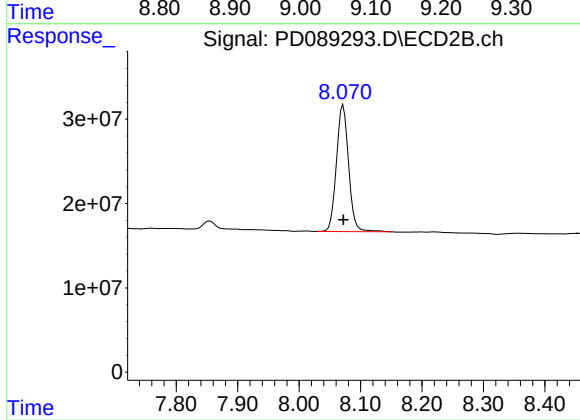
#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.000 min
Response: 41602422
Conc: 10.59 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-66

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 206274214
Conc: 10.43 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-60	SDG No.:	Q2458
Lab Sample ID:	Q2458-06	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	92.5 Decanted:
Sample Wt/Vol:	30.05 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089294.D	1	07/01/25 08:30	07/01/25 21:10	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.80	ug/kg
319-85-7	beta-BHC	0.19	U	0.19	1.80	ug/kg
319-86-8	delta-BHC	0.42	U	0.42	1.80	ug/kg
58-89-9	gamma-BHC (Lindane)	0.15	U	0.15	1.80	ug/kg
76-44-8	Heptachlor	0.13	U	0.13	1.80	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.80	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.80	ug/kg
959-98-8	Endosulfan I	0.15	U	0.15	1.80	ug/kg
60-57-1	Dieldrin	0.15	U	0.15	1.80	ug/kg
72-55-9	4,4-DDE	0.15	U	0.15	1.80	ug/kg
72-20-8	Endrin	0.15	U	0.15	1.80	ug/kg
33213-65-9	Endosulfan II	0.31	U	0.31	1.80	ug/kg
72-54-8	4,4-DDD	0.17	J	0.16	1.80	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.80	ug/kg
50-29-3	4,4-DDT	0.36	JP	0.15	1.80	ug/kg
72-43-5	Methoxychlor	0.40	U	0.40	1.80	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.80	ug/kg
7421-93-4	Endrin aldehyde	0.40	U	0.40	1.80	ug/kg
5103-71-9	alpha-Chlordane	0.90	JP	0.13	1.80	ug/kg
5103-74-2	gamma-Chlordane	0.46	J	0.16	1.80	ug/kg
8001-35-2	Toxaphene	5.80	U	5.80	35.6	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	11.6		20 - 144	58%	SPK: 20
877-09-8	Tetrachloro-m-xylene	13.7		19 - 148	68%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-60	SDG No.:	Q2458
Lab Sample ID:	Q2458-06	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	92.5 Decanted:
Sample Wt/Vol:	30.05 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089294.D	1	07/01/25 08:30	07/01/25 21:10	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089294.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 21:10
 Operator : AR\AJ
 Sample : Q2458-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-60

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/02/2025
 Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:52:12 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.550	2.881	35343982	231.5E6	12.388	13.653
2) SA Decachlor...	9.073	8.072	45383904	146.8E6	11.555	7.426 #
Target Compounds						
10) B gamma-Chl...	5.944	5.124	6128061	21242718	1.280m	0.888m#
11) B alpha-Chl...	6.030	5.188	12141915	36822266	2.509m	1.608m#
16) A 4,4'-DDD	6.700	5.927	1350255	8828695	0.393m	0.462m
17) MA 4,4'-DDT	7.019	6.180	1592517	20054610	0.420m	0.989 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089294.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 21:10
Operator : AR\AJ
Sample : Q2458-06
Misc :
ALS Vial : 26 Sample Multiplier: 1

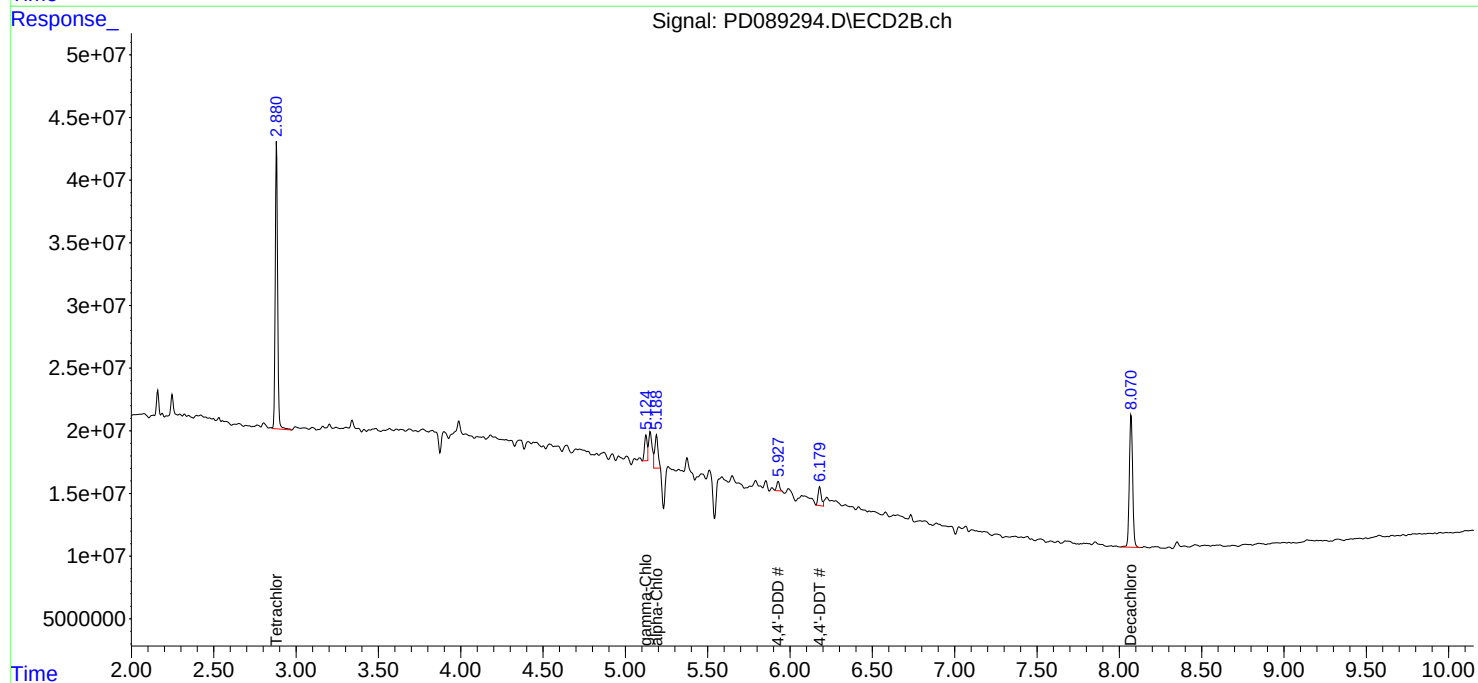
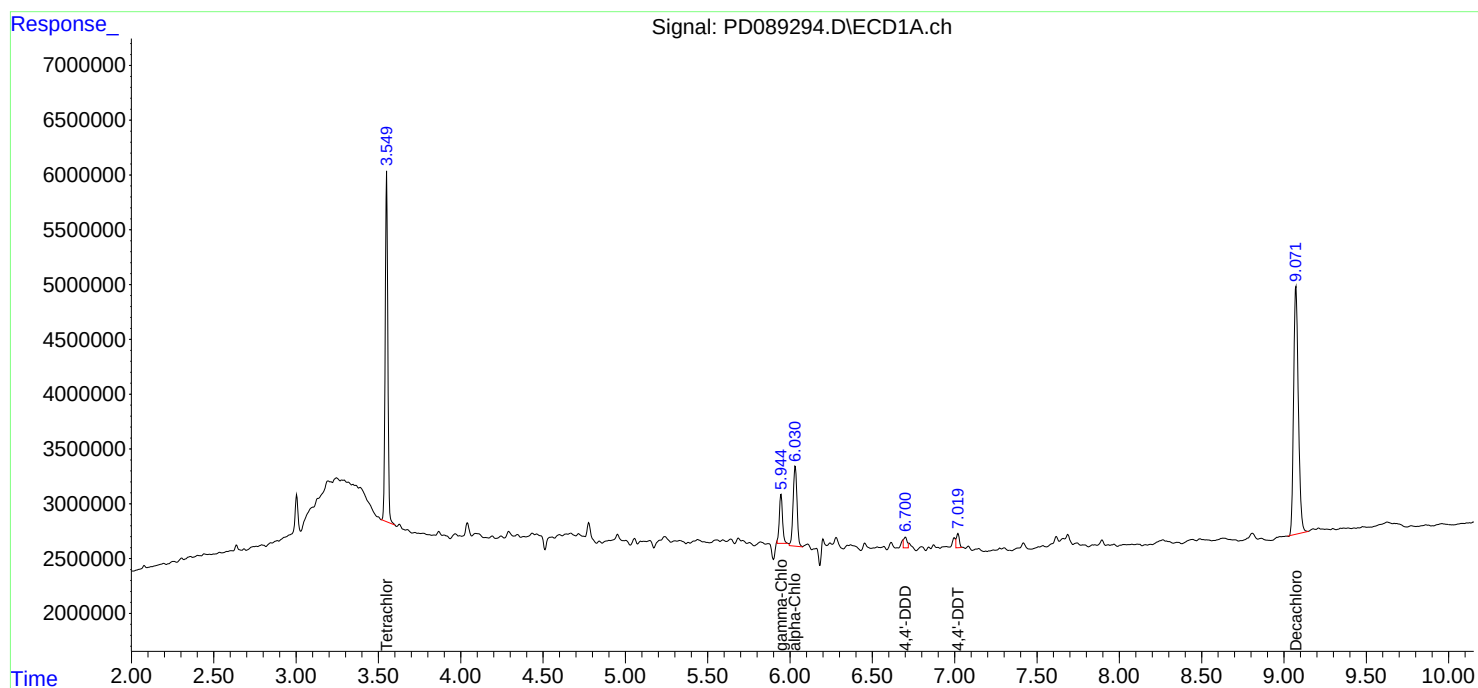
Instrument :
ECD_D
ClientSampleId :
TP-60

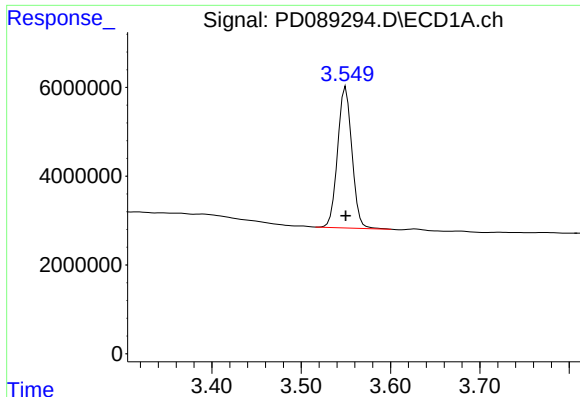
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:12 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





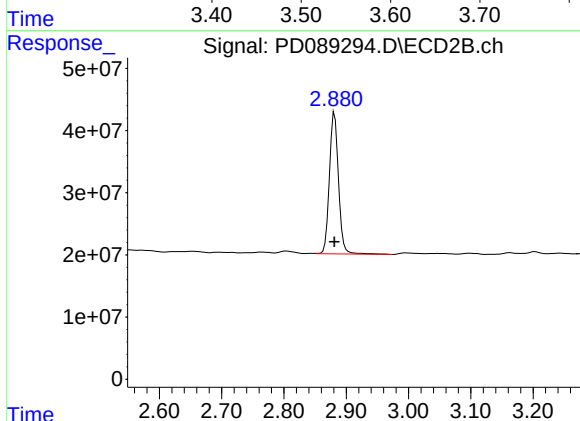
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 35343982
Conc: 12.39 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-60

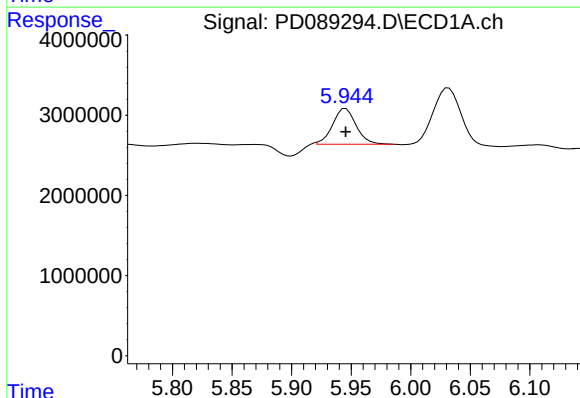
Manual Integrations
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Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



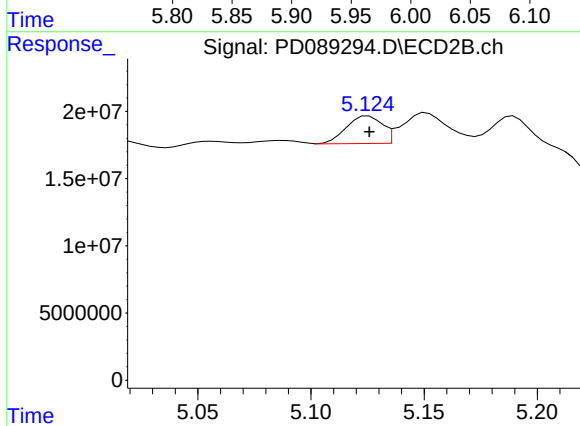
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 231465961
Conc: 13.65 ng/ml



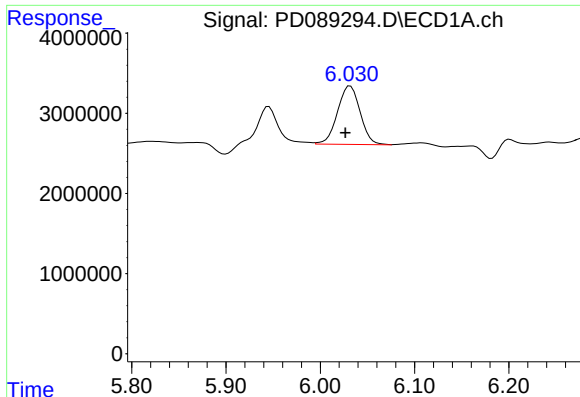
#10 gamma-Chlordane

R.T.: 5.944 min
Delta R.T.: -0.001 min
Response: 6128061
Conc: 1.28 ng/ml m



#10 gamma-Chlordane

R.T.: 5.124 min
Delta R.T.: -0.002 min
Response: 21242718
Conc: 0.89 ng/ml m



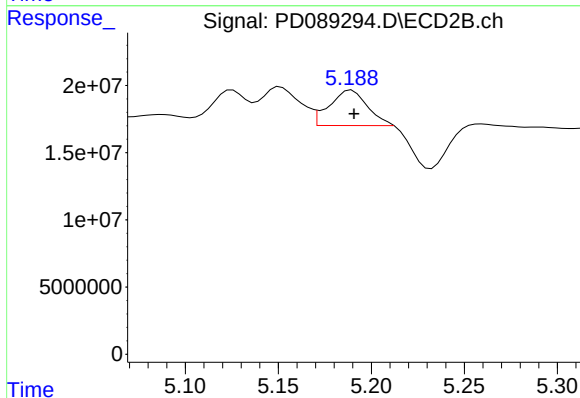
#11 alpha-Chlordane

R.T.: 6.030 min
Delta R.T.: 0.004 min
Response: 12141915
Conc: 2.51 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-60

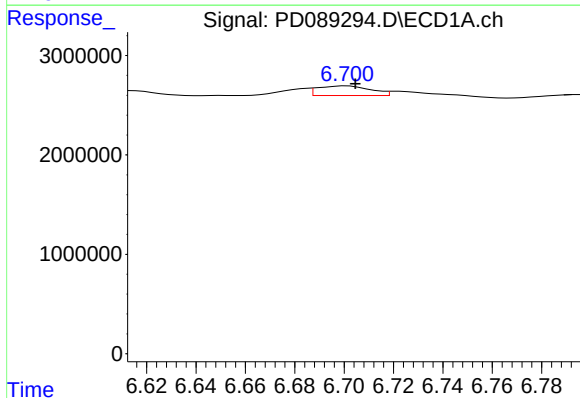
Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



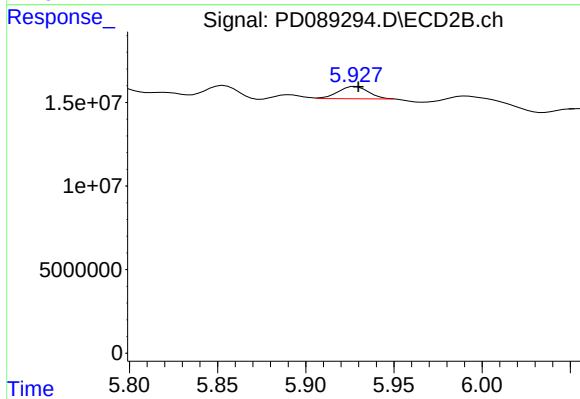
#11 alpha-Chlordane

R.T.: 5.188 min
Delta R.T.: -0.003 min
Response: 36822266
Conc: 1.61 ng/ml m



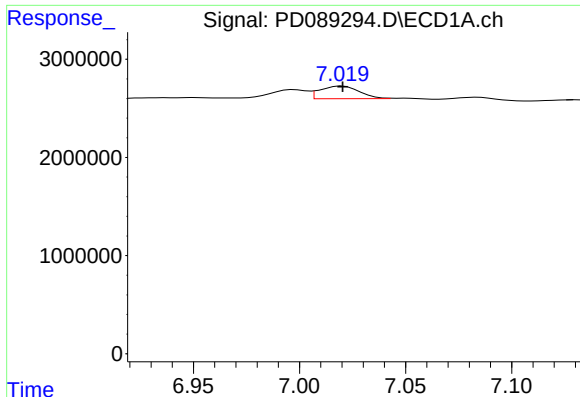
#16 4,4'-DDD

R.T.: 6.700 min
Delta R.T.: -0.005 min
Response: 1350255
Conc: 0.39 ng/ml m



#16 4,4'-DDD

R.T.: 5.927 min
Delta R.T.: -0.003 min
Response: 8828695
Conc: 0.46 ng/ml m



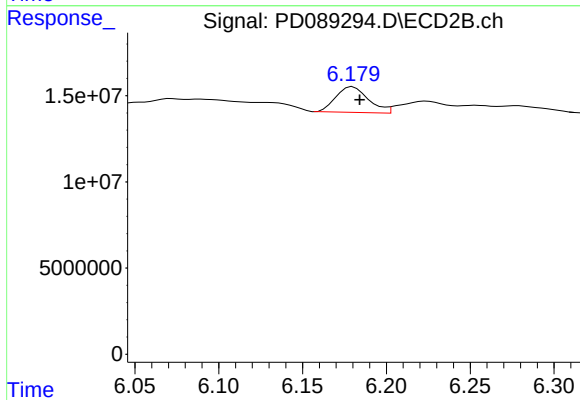
#17 4,4'-DDT

R.T.: 7.019 min
Delta R.T.: -0.002 min
Response: 1592517
Conc: 0.42 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-60

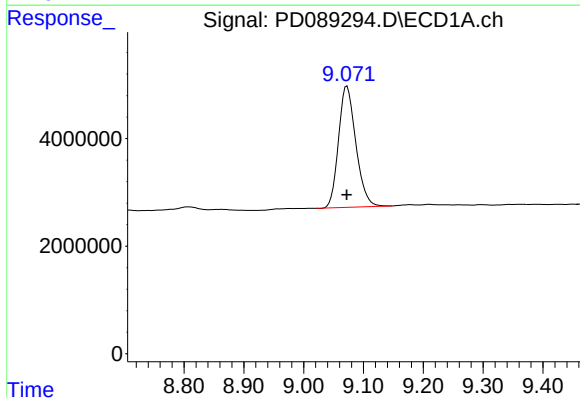
Manual Integrations
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Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



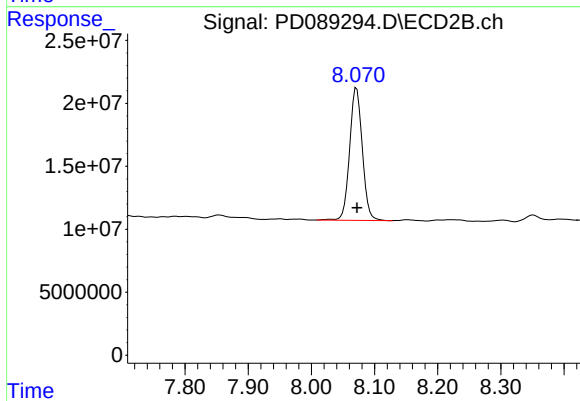
#17 4,4'-DDT

R.T.: 6.180 min
Delta R.T.: -0.004 min
Response: 20054610
Conc: 0.99 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.000 min
Response: 45383904
Conc: 11.56 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 146805452
Conc: 7.43 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-62	SDG No.:	Q2458
Lab Sample ID:	Q2458-07	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	91.1 Decanted:
Sample Wt/Vol:	30.04 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089295.D	1	07/01/25 08:30	07/01/25 21:24	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.14	U	0.14	1.90	ug/kg
319-85-7	beta-BHC	0.20	U	0.20	1.90	ug/kg
319-86-8	delta-BHC	0.43	U	0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	0.15	U	0.15	1.90	ug/kg
76-44-8	Heptachlor	0.17	J	0.13	1.90	ug/kg
309-00-2	Aldrin	0.13	U	0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	0.21	U	0.21	1.90	ug/kg
959-98-8	Endosulfan I	0.15	U	0.15	1.90	ug/kg
60-57-1	Dieldrin	0.15	U	0.15	1.90	ug/kg
72-55-9	4,4-DDE	0.15	U	0.15	1.90	ug/kg
72-20-8	Endrin	0.15	U	0.15	1.90	ug/kg
33213-65-9	Endosulfan II	0.32	U	0.32	1.90	ug/kg
72-54-8	4,4-DDD	0.16	U	0.16	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	0.14	U	0.14	1.90	ug/kg
50-29-3	4,4-DDT	0.15	U	0.15	1.90	ug/kg
72-43-5	Methoxychlor	0.41	U	0.41	1.90	ug/kg
53494-70-5	Endrin ketone	0.21	U	0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	0.41	U	0.41	1.90	ug/kg
5103-71-9	alpha-Chlordane	0.13	U	0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	0.16	U	0.16	1.90	ug/kg
8001-35-2	Toxaphene	5.90	U	5.90	36.2	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	12.5		20 - 144	62%	SPK: 20
877-09-8	Tetrachloro-m-xylene	17.6		19 - 148	88%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-62	SDG No.:	Q2458
Lab Sample ID:	Q2458-07	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	91.1 Decanted:
Sample Wt/Vol:	30.04 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089295.D	1	07/01/25 08:30	07/01/25 21:24	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089295.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 21:24
 Operator : AR\AJ
 Sample : Q2458-07
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-62

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:52:24 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.550	2.880	44933438	298.9E6	15.749	17.630m
28) SA Decachlor...	9.072	8.071	48993057	207.4E6	12.474	10.491
Target Compounds						
4) MA Heptachlor	4.936	4.082	2492245	8614864	0.457	0.344

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089295.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 21:24
Operator : AR\AJ
Sample : Q2458-07
Misc :
ALS Vial : 27 Sample Multiplier: 1

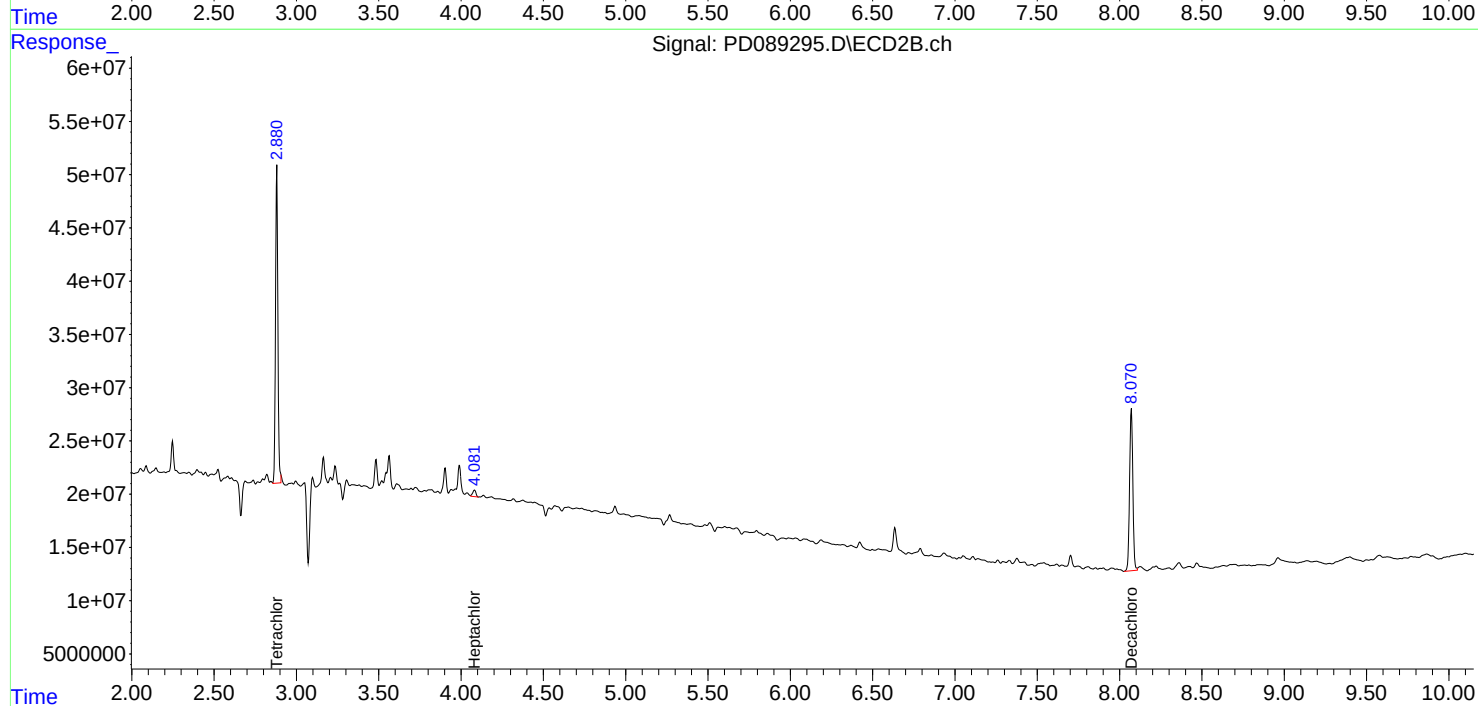
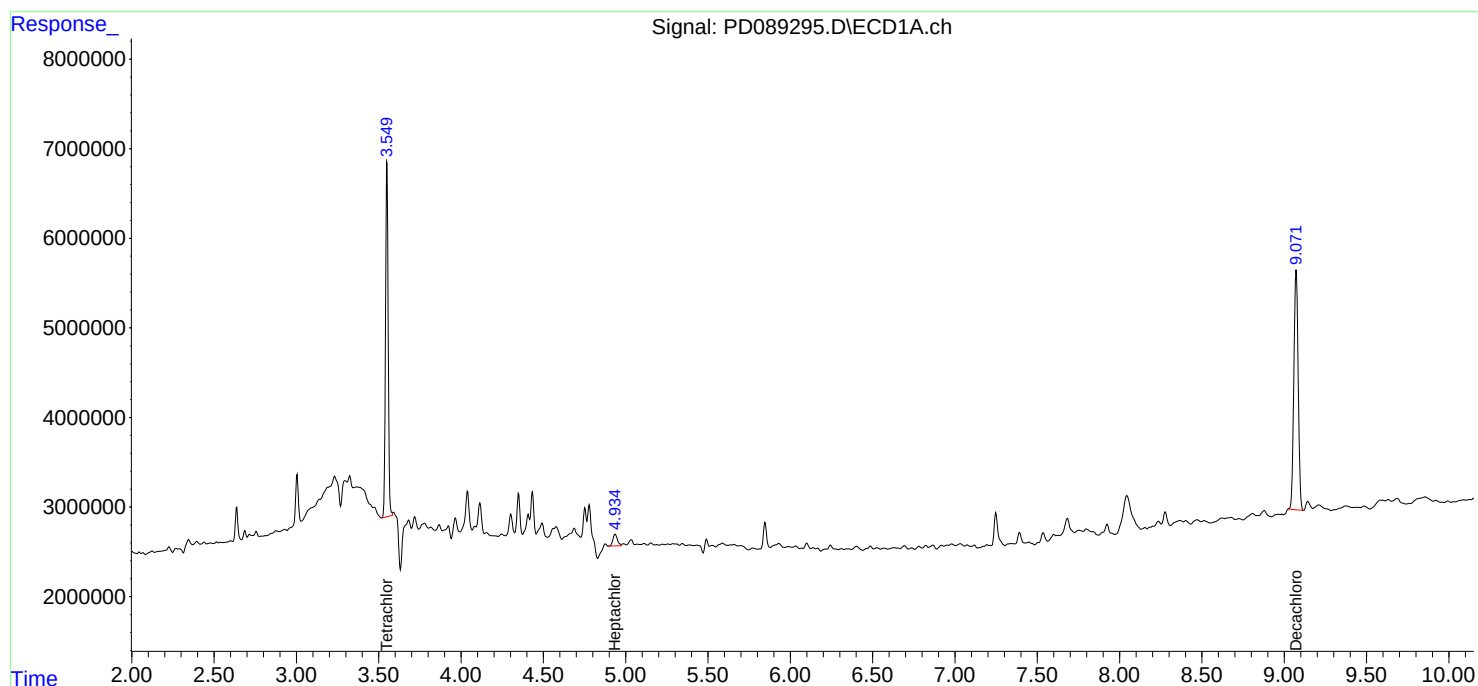
Instrument :
ECD_D
ClientSampleId :
TP-62

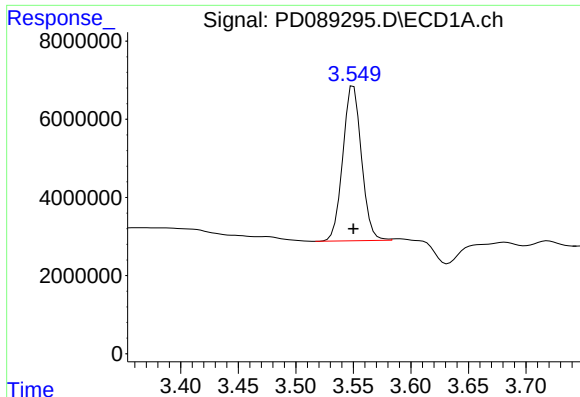
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:24 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





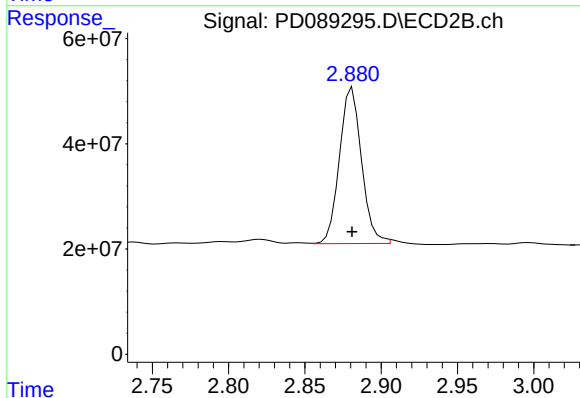
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
 Delta R.T.: 0.000 min
 Response: 44933438
 Conc: 15.75 ng/ml

Instrument :
 ECD_D
 ClientSampleId :
 TP-62

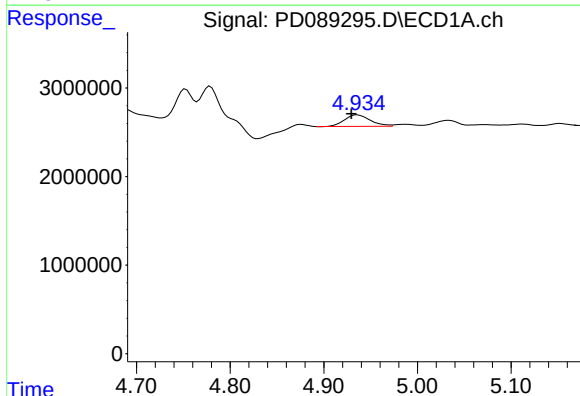
Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025



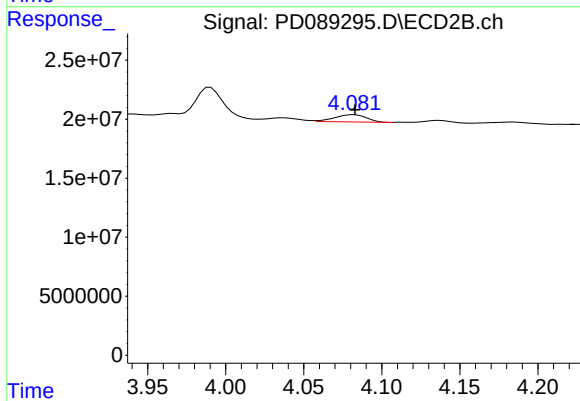
#1 Tetrachloro-m-xylene

R.T.: 2.880 min
 Delta R.T.: -0.001 min
 Response: 298878304
 Conc: 17.63 ng/ml m



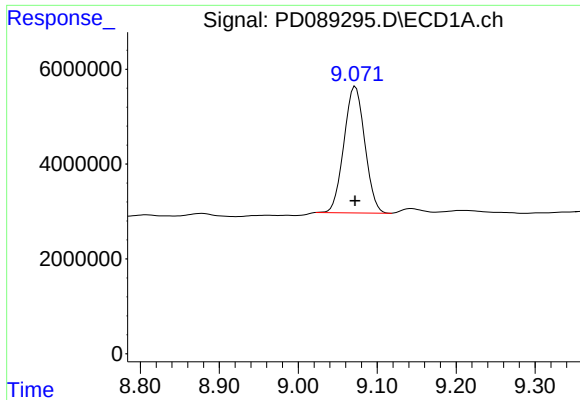
#4 Heptachlor

R.T.: 4.936 min
 Delta R.T.: 0.006 min
 Response: 2492245
 Conc: 0.46 ng/ml



#4 Heptachlor

R.T.: 4.082 min
 Delta R.T.: 0.000 min
 Response: 8614864
 Conc: 0.34 ng/ml



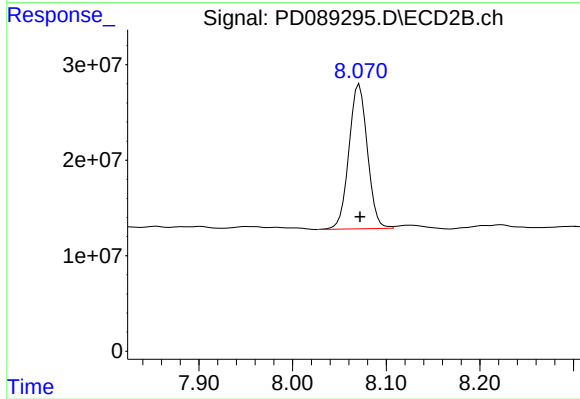
#28 Decachlorobiphenyl

R.T.: 9.072 min
Delta R.T.: 0.000 min
Response: 48993057
Conc: 12.47 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-62

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



#28 Decachlorobiphenyl

R.T.: 8.071 min
Delta R.T.: 0.000 min
Response: 207402121
Conc: 10.49 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-63	SDG No.:	Q2458
Lab Sample ID:	Q2458-08	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	86.4 Decanted:
Sample Wt/Vol:	30.08 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089296.D	1	07/01/25 08:30	07/01/25 21:37	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.15	U	0.15	2.00	ug/kg
319-85-7	beta-BHC	0.21	U	0.21	2.00	ug/kg
319-86-8	delta-BHC	0.45	U	0.45	2.00	ug/kg
58-89-9	gamma-BHC (Lindane)	0.16	U	0.16	2.00	ug/kg
76-44-8	Heptachlor	0.14	U	0.14	2.00	ug/kg
309-00-2	Aldrin	0.14	U	0.14	2.00	ug/kg
1024-57-3	Heptachlor epoxide	0.22	U	0.22	2.00	ug/kg
959-98-8	Endosulfan I	0.16	U	0.16	2.00	ug/kg
60-57-1	Dieldrin	0.16	U	0.16	2.00	ug/kg
72-55-9	4,4-DDE	0.16	U	0.16	2.00	ug/kg
72-20-8	Endrin	0.16	U	0.16	2.00	ug/kg
33213-65-9	Endosulfan II	0.33	U	0.33	2.00	ug/kg
72-54-8	4,4-DDD	0.41	J	0.17	2.00	ug/kg
1031-07-8	Endosulfan Sulfate	0.15	U	0.15	2.00	ug/kg
50-29-3	4,4-DDT	0.44	JP	0.16	2.00	ug/kg
72-43-5	Methoxychlor	0.43	U	0.43	2.00	ug/kg
53494-70-5	Endrin ketone	0.22	U	0.22	2.00	ug/kg
7421-93-4	Endrin aldehyde	0.43	U	0.43	2.00	ug/kg
5103-71-9	alpha-Chlordane	0.58	J	0.14	2.00	ug/kg
5103-74-2	gamma-Chlordane	0.40	J	0.17	2.00	ug/kg
8001-35-2	Toxaphene	6.20	U	6.20	38.1	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	10.2		20 - 144	51%	SPK: 20
877-09-8	Tetrachloro-m-xylene	12.7		19 - 148	64%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-63	SDG No.:	Q2458
Lab Sample ID:	Q2458-08	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	86.4 Decanted:
Sample Wt/Vol:	30.08 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089296.D	1	07/01/25 08:30	07/01/25 21:37	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
------------	-----------	-------	-----------	-----	------------	-------

Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089296.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 21:37
 Operator : AR\AJ
 Sample : Q2458-08
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-63

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/02/2025
 Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:52:35 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.880	33016136	216.0E6	11.572	12.740m
28)	SA Decachlor...	9.073	8.071	39895212	151.6E6	10.158	7.668
Target Compounds							
10)	B gamma-Chl...	5.944	5.124	4921330	18057708	1.028m	0.755m#
11)	B alpha-Chl...	6.029	5.188	7357490	27737576	1.521	1.211m
14)	MA Endrin	6.574	5.793	1364116	6827957	0.331	0.314
16)	A 4,4'-DDD	6.703	5.929	3682947	15945348	1.073m	0.834
17)	MA 4,4'-DDT	7.020	6.180	2388886	23188038	0.631	1.143 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089296.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 21:37
Operator : AR\AJ
Sample : Q2458-08
Misc :
ALS Vial : 28 Sample Multiplier: 1

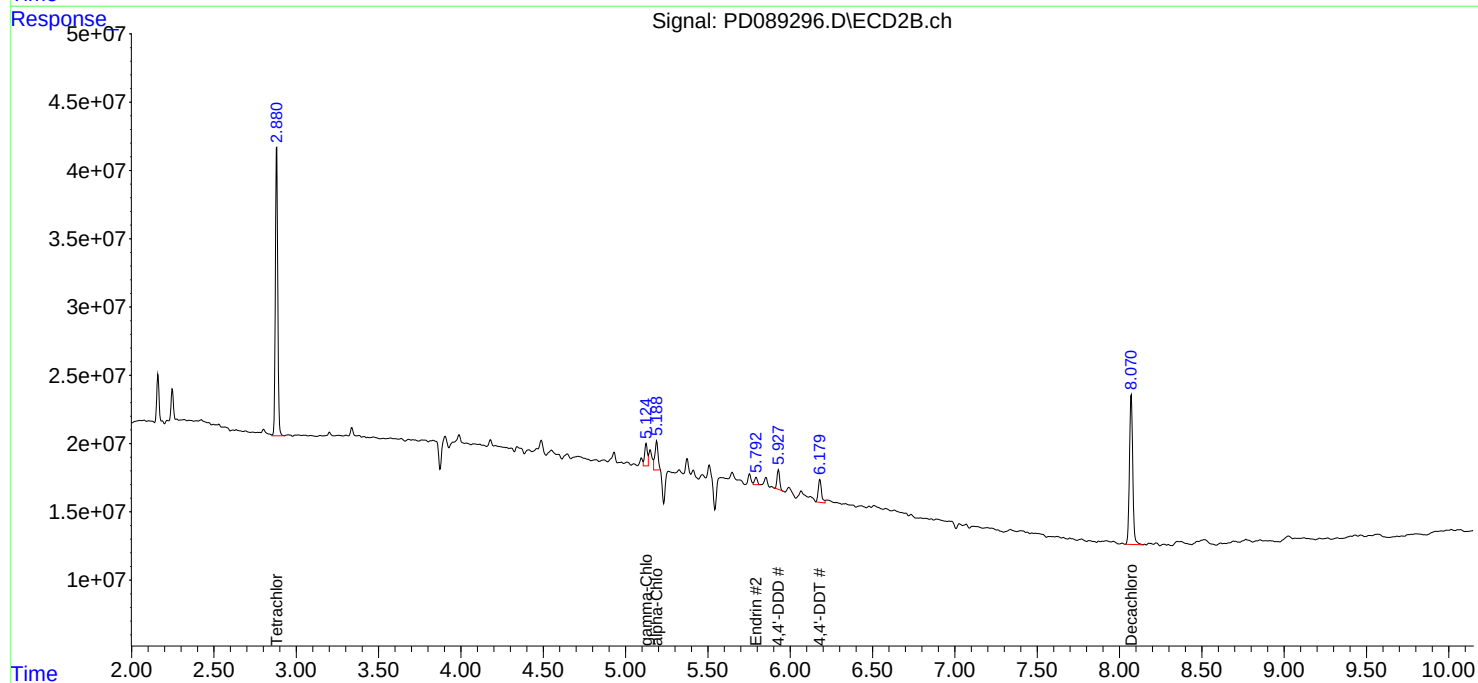
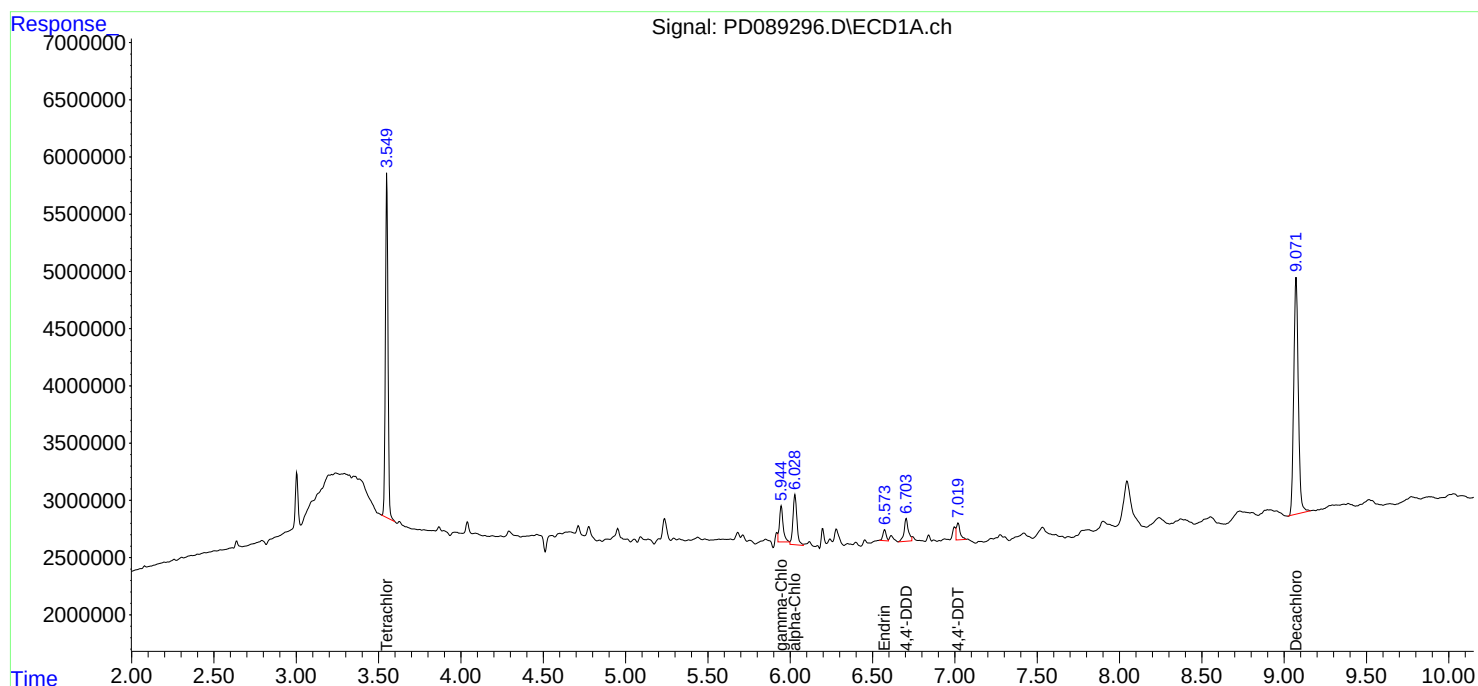
Instrument :
ECD_D
ClientSampleId :
TP-63

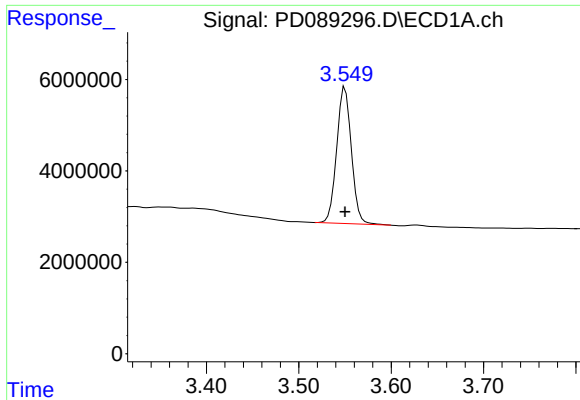
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:35 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





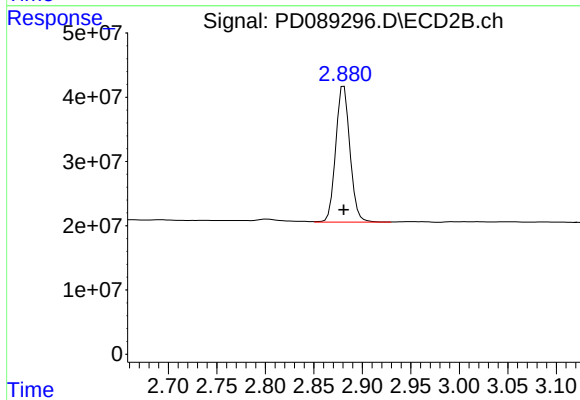
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 33016136
Conc: 11.57 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-63

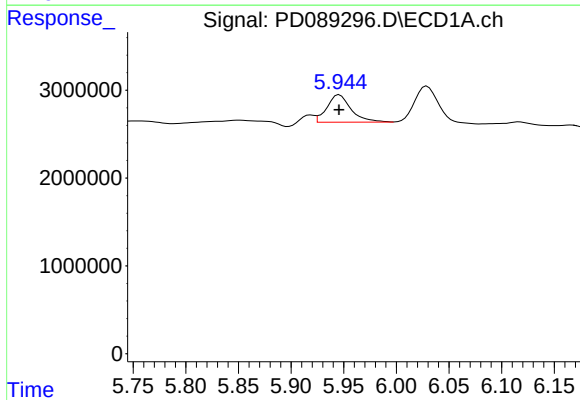
Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



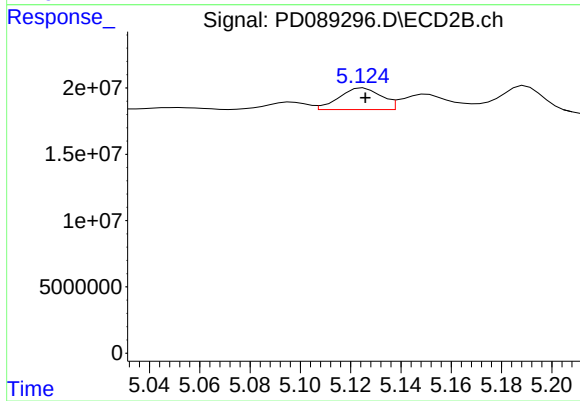
#1 Tetrachloro-m-xylene

R.T.: 2.880 min
Delta R.T.: -0.001 min
Response: 215977192
Conc: 12.74 ng/ml m



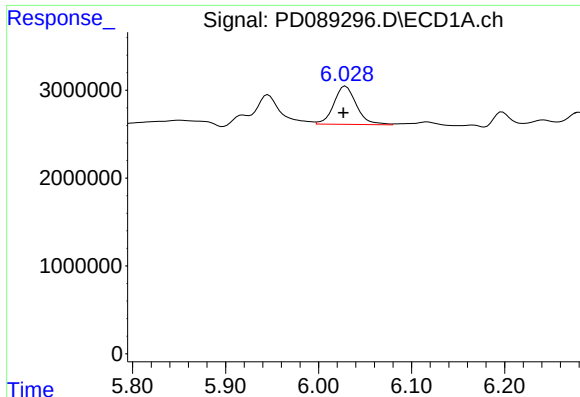
#10 gamma-Chlordane

R.T.: 5.944 min
Delta R.T.: -0.001 min
Response: 4921330
Conc: 1.03 ng/ml m



#10 gamma-Chlordane

R.T.: 5.124 min
Delta R.T.: -0.002 min
Response: 18057708
Conc: 0.75 ng/ml m



#11 alpha-Chlordane

R.T.: 6.029 min
Delta R.T.: 0.002 min
Response: 7357490
Conc: 1.52 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-63

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025

#11 alpha-Chlordane

R.T.: 5.188 min
Delta R.T.: -0.003 min
Response: 27737576
Conc: 1.21 ng/ml m

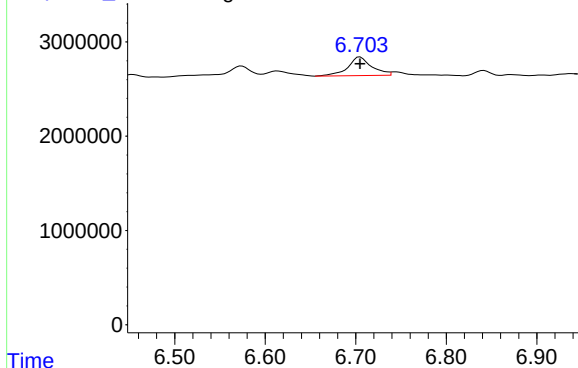
#14 Endrin

R.T.: 6.574 min
Delta R.T.: 0.000 min
Response: 1364116
Conc: 0.33 ng/ml

#14 Endrin

R.T.: 5.793 min
Delta R.T.: 0.004 min
Response: 6827957
Conc: 0.31 ng/ml

Response_ Signal: PD089296.D\ECD1A.ch



#16 4,4'-DDD

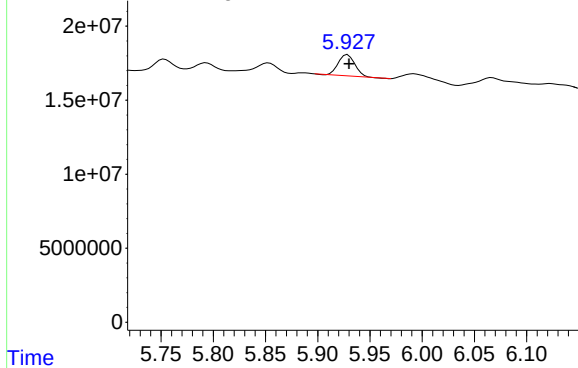
R.T.: 6.703 min
Delta R.T.: -0.001 min
Response: 3682947
Conc: 1.07 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-63

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025

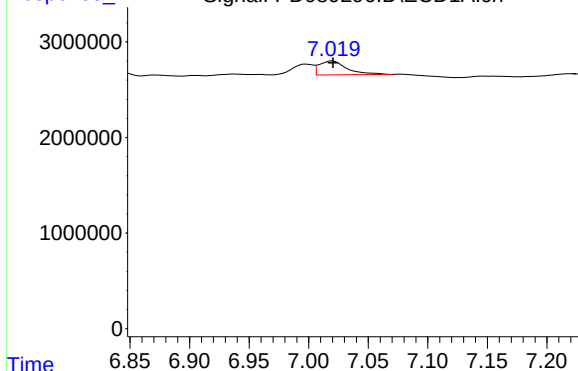
Response_ Signal: PD089296.D\ECD2B.ch



#16 4,4'-DDD

R.T.: 5.929 min
Delta R.T.: -0.001 min
Response: 15945348
Conc: 0.83 ng/ml

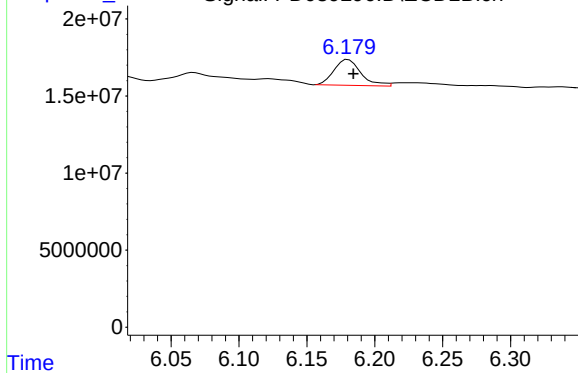
Response_ Signal: PD089296.D\ECD1A.ch



#17 4,4'-DDT

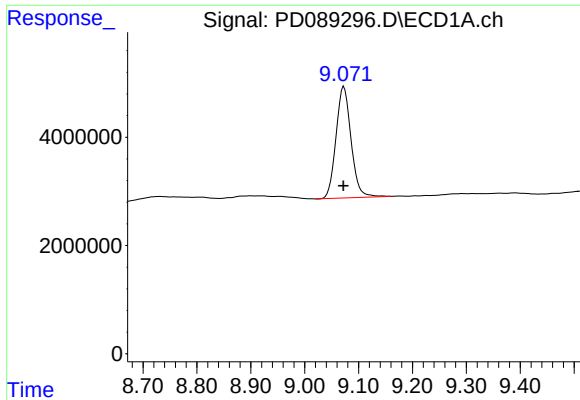
R.T.: 7.020 min
Delta R.T.: 0.000 min
Response: 2388886
Conc: 0.63 ng/ml

Response_ Signal: PD089296.D\ECD2B.ch



#17 4,4'-DDT

R.T.: 6.180 min
Delta R.T.: -0.004 min
Response: 23188038
Conc: 1.14 ng/ml



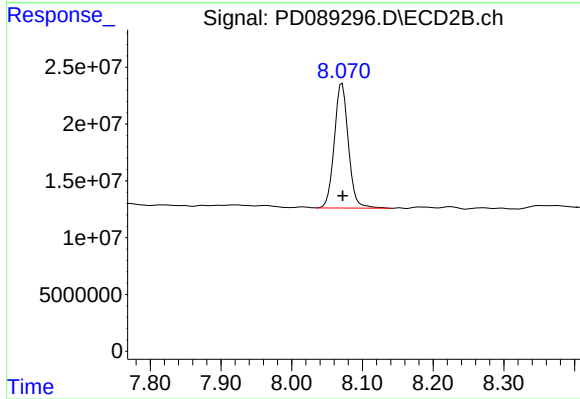
#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.000 min
Response: 39895212
Conc: 10.16 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-63

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



#28 Decachlorobiphenyl

R.T.: 8.071 min
Delta R.T.: 0.000 min
Response: 151599662
Conc: 7.67 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-59	SDG No.:	Q2458
Lab Sample ID:	Q2458-09	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	76.6 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089297.D	1	07/01/25 08:30	07/01/25 21:51	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.17	U	0.17	2.20	ug/kg
319-85-7	beta-BHC	0.23	U	0.23	2.20	ug/kg
319-86-8	delta-BHC	0.51	U	0.51	2.20	ug/kg
58-89-9	gamma-BHC (Lindane)	0.18	U	0.18	2.20	ug/kg
76-44-8	Heptachlor	0.16	U	0.16	2.20	ug/kg
309-00-2	Aldrin	0.16	U	0.16	2.20	ug/kg
1024-57-3	Heptachlor epoxide	0.25	U	0.25	2.20	ug/kg
959-98-8	Endosulfan I	0.18	U	0.18	2.20	ug/kg
60-57-1	Dieldrin	0.18	U	0.18	2.20	ug/kg
72-55-9	4,4-DDE	0.18	U	0.18	2.20	ug/kg
72-20-8	Endrin	0.18	U	0.18	2.20	ug/kg
33213-65-9	Endosulfan II	0.38	U	0.38	2.20	ug/kg
72-54-8	4,4-DDD	0.20	U	0.20	2.20	ug/kg
1031-07-8	Endosulfan Sulfate	0.17	U	0.17	2.20	ug/kg
50-29-3	4,4-DDT	0.18	U	0.18	2.20	ug/kg
72-43-5	Methoxychlor	0.48	U	0.48	2.20	ug/kg
53494-70-5	Endrin ketone	0.25	U	0.25	2.20	ug/kg
7421-93-4	Endrin aldehyde	0.48	U	0.48	2.20	ug/kg
5103-71-9	alpha-Chlordane	0.16	U	0.16	2.20	ug/kg
5103-74-2	gamma-Chlordane	0.20	U	0.20	2.20	ug/kg
8001-35-2	Toxaphene	7.10	U	7.10	43.0	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	12.9		20 - 144	65%	SPK: 20
877-09-8	Tetrachloro-m-xylene	19.0		19 - 148	95%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-59	SDG No.:	Q2458
Lab Sample ID:	Q2458-09	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	76.6 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089297.D	1	07/01/25 08:30	07/01/25 21:51	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089297.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 21:51
 Operator : AR\AJ
 Sample : Q2458-09
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-59

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:52:47 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.879	50314076	321.4E6	17.635	18.956m
28)	SA Decachlor...	9.073	8.072	50856481	225.0E6	12.948	11.383

Target Compounds

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089297.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 21:51
Operator : AR\AJ
Sample : Q2458-09
Misc :
ALS Vial : 29 Sample Multiplier: 1

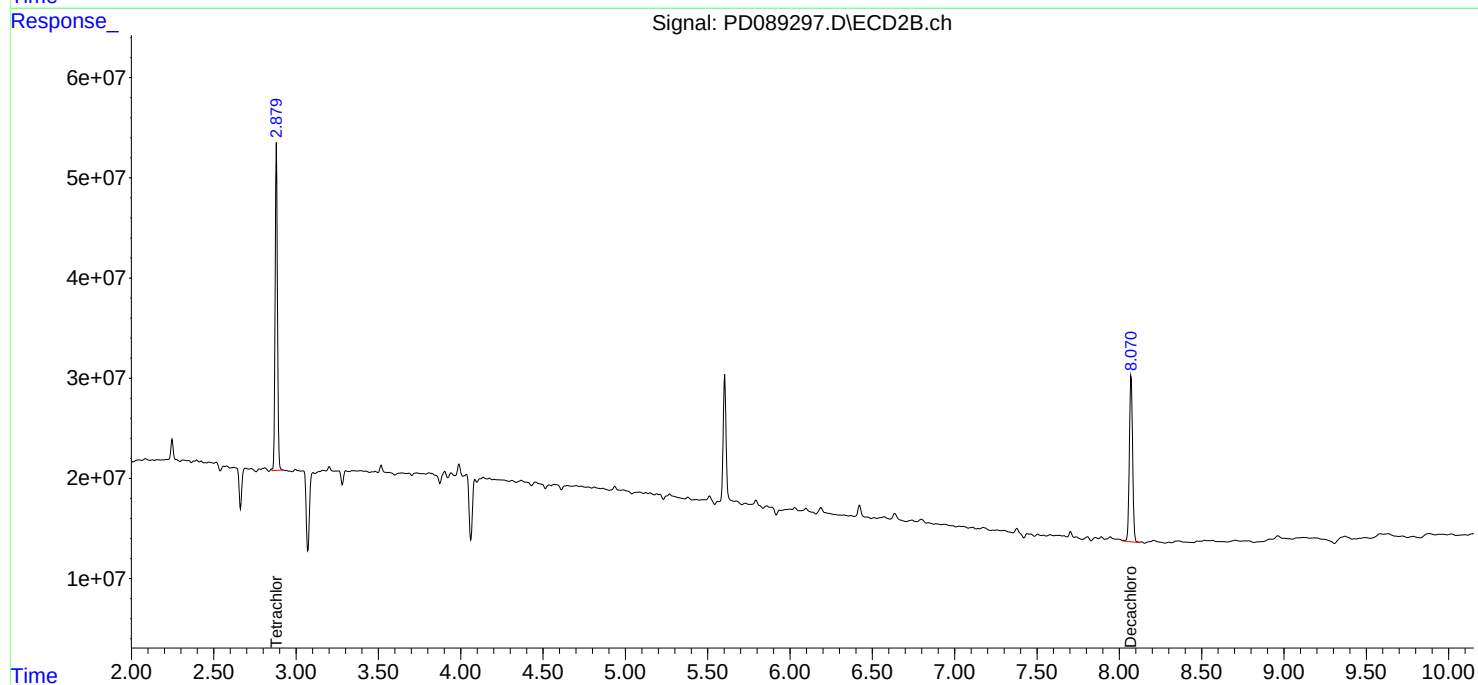
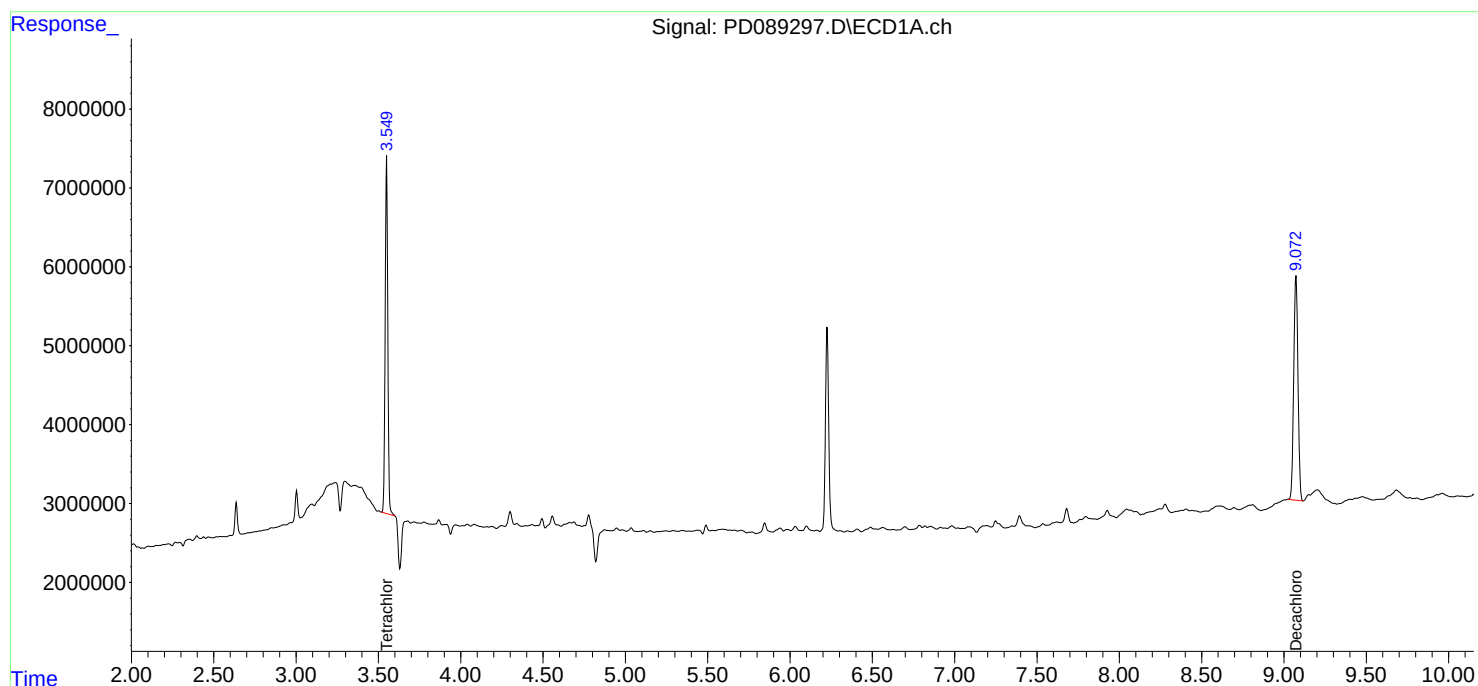
Instrument :
ECD_D
ClientSampleId :
TP-59

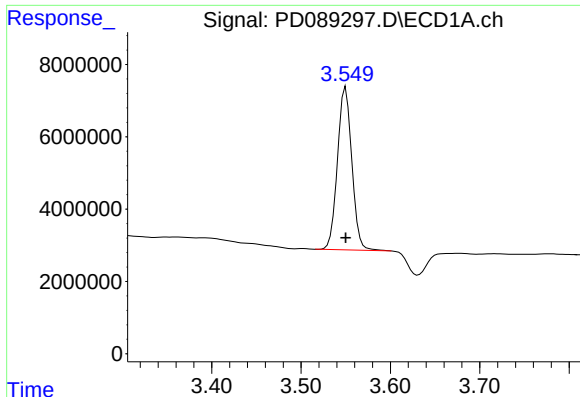
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:47 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





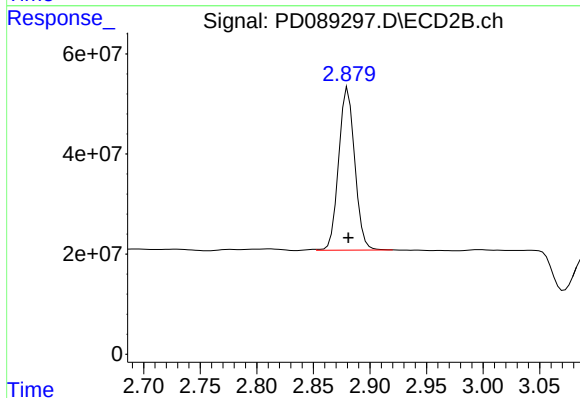
#1 Tetrachloro-m-xylene

R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 50314076
Conc: 17.63 ng/ml

Instrument :
ECD_D
ClientSampleId :
TP-59

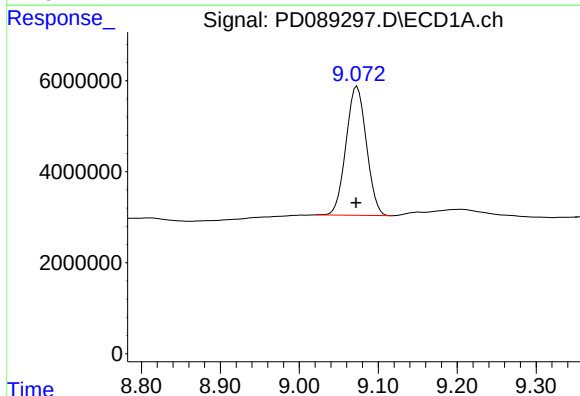
Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/02/2025
Supervised By :mohammad ahmed 07/04/2025



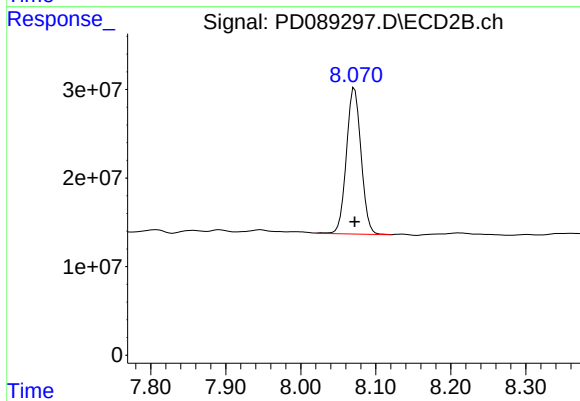
#1 Tetrachloro-m-xylene

R.T.: 2.879 min
Delta R.T.: -0.002 min
Response: 321361786
Conc: 18.96 ng/ml m



#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.001 min
Response: 50856481
Conc: 12.95 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 225039940
Conc: 11.38 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	FB-06272025	SDG No.:	Q2458
Lab Sample ID:	Q2458-10	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	980 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089332.D	1	07/03/25 08:58	07/03/25 14:40	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0040	U	0.0040	0.051	ug/L
319-85-7	beta-BHC	0.0050	U	0.0050	0.051	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.051	ug/L
58-89-9	gamma-BHC (Lindane)	0.0038	U	0.0038	0.051	ug/L
76-44-8	Heptachlor	0.0028	U	0.0028	0.051	ug/L
309-00-2	Aldrin	0.0037	U	0.0037	0.051	ug/L
1024-57-3	Heptachlor epoxide	0.0098	U	0.0098	0.051	ug/L
959-98-8	Endosulfan I	0.0032	U	0.0032	0.051	ug/L
60-57-1	Dieldrin	0.0037	U	0.0037	0.051	ug/L
72-55-9	4,4-DDE	0.0038	U	0.0038	0.051	ug/L
72-20-8	Endrin	0.0033	U	0.0033	0.051	ug/L
33213-65-9	Endosulfan II	0.0081	U	0.0081	0.051	ug/L
72-54-8	4,4-DDD	0.0072	U	0.0072	0.051	ug/L
1031-07-8	Endosulfan Sulfate	0.0038	U	0.0038	0.051	ug/L
50-29-3	4,4-DDT	0.0036	U	0.0036	0.051	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.051	ug/L
53494-70-5	Endrin ketone	0.0095	U	0.0095	0.051	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.051	ug/L
5103-71-9	alpha-Chlordane	0.0036	U	0.0036	0.051	ug/L
5103-74-2	gamma-Chlordane	0.0040	U	0.0040	0.051	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	13.7		57 - 171	69%	SPK: 20
877-09-8	Tetrachloro-m-xylene	21.0		61 - 148	105%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	FB-06272025	SDG No.:	Q2458
Lab Sample ID:	Q2458-10	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	980	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089332.D	1	07/03/25 08:58	07/03/25 14:40	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
------------	-----------	-------	-----------	-----	------------	-------

Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089332.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 14:40
 Operator : AR\AJ
 Sample : Q2458-10
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 FB-06272025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:18:12 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.881	52768868	356.4E6	18.495	21.022
28)	SA Decachlor...	9.075	8.072	51727903	270.9E6	13.170	13.700

Target Compounds

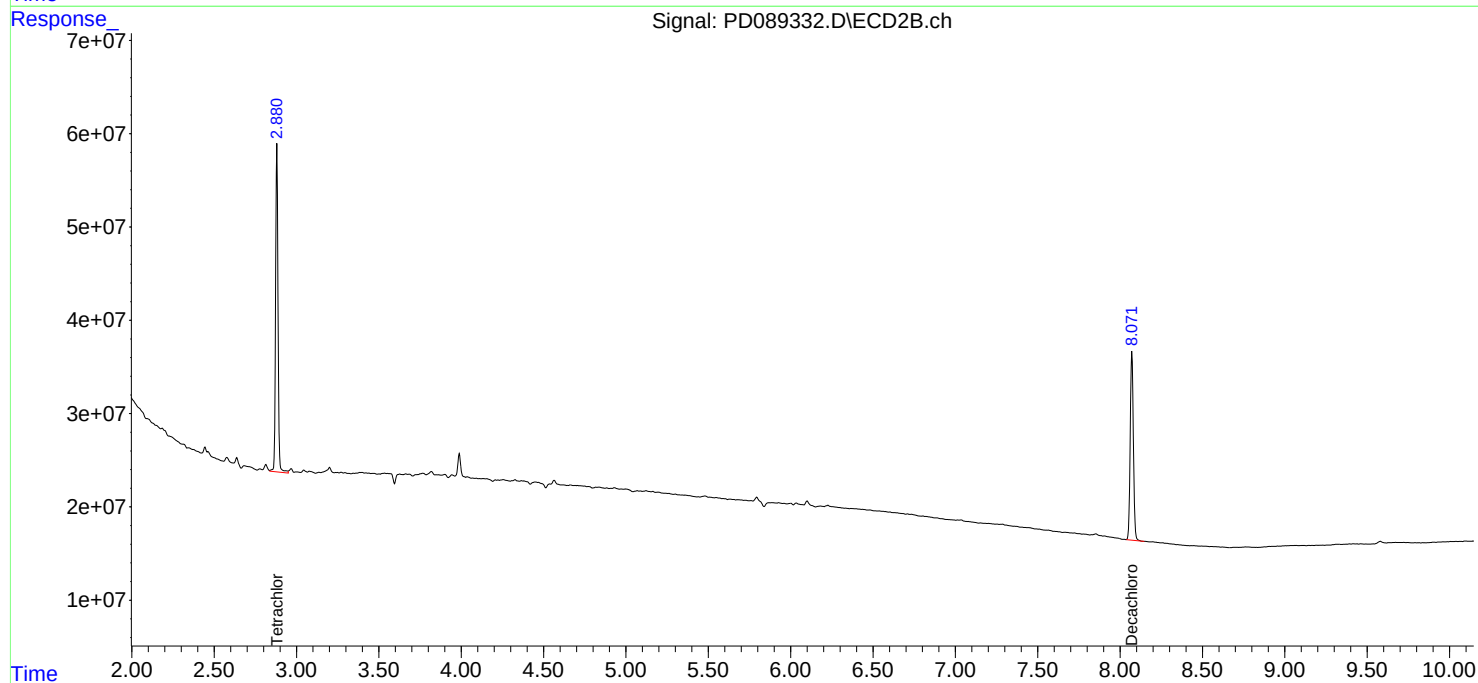
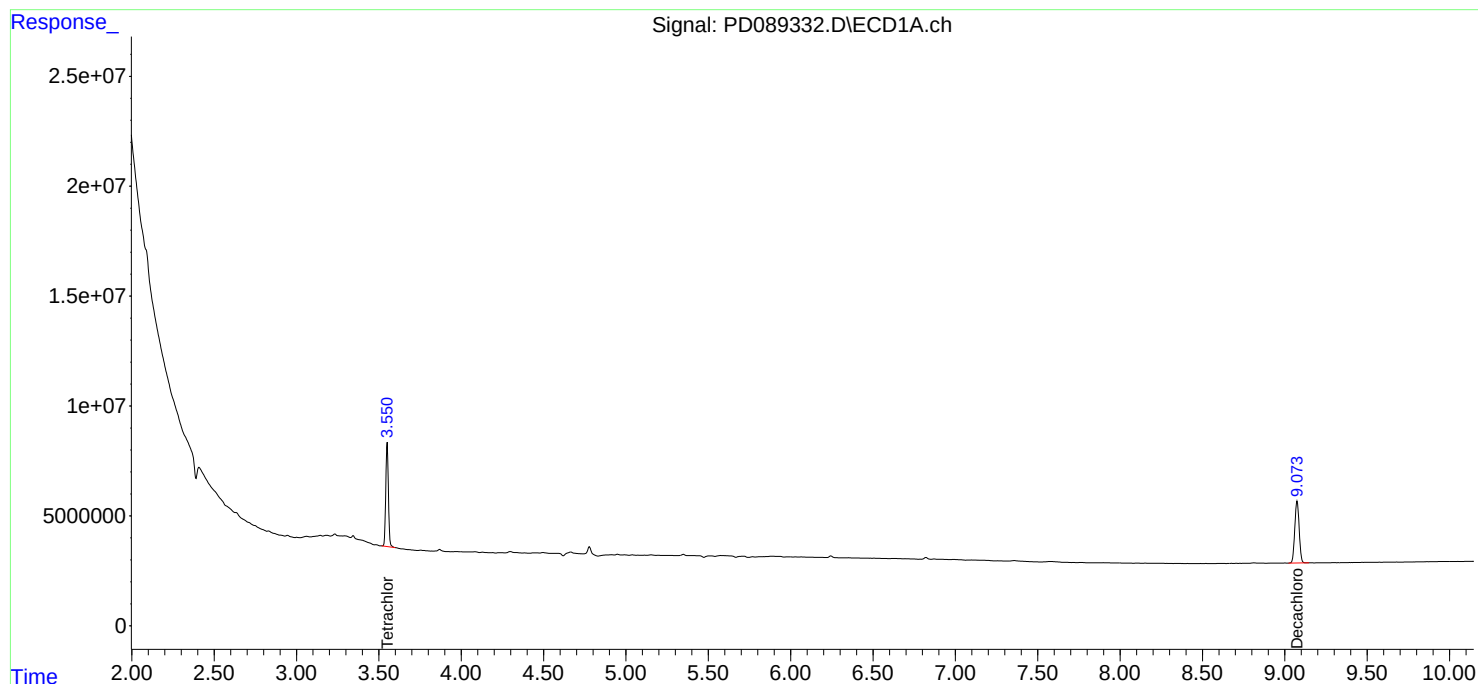
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

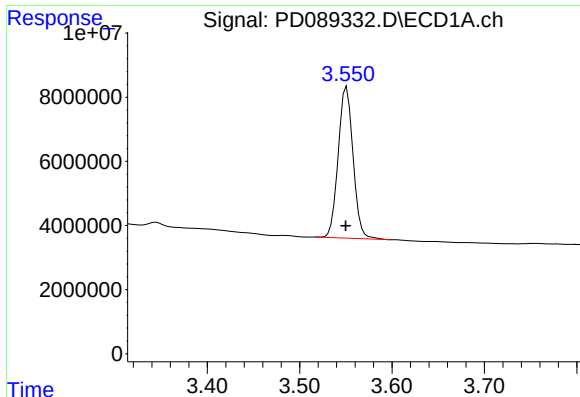
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089332.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 14:40
Operator : AR\AJ
Sample : Q2458-10
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
FB-06272025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:18:12 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

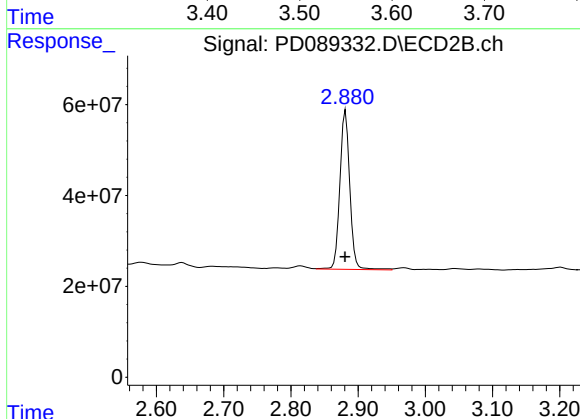




#1 Tetrachloro-m-xylene

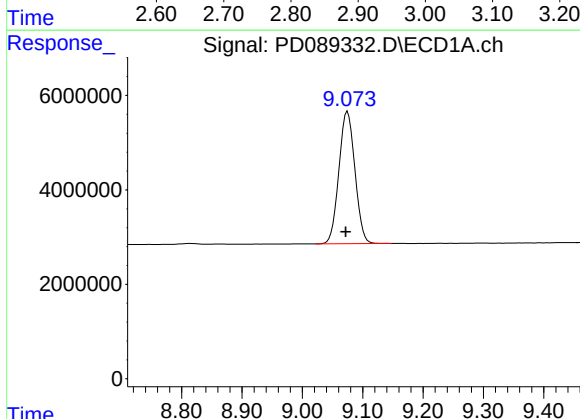
R.T.: 3.551 min
Delta R.T.: 0.000 min
Response: 52768868
Conc: 18.50 ng/ml

Instrument :
ECD_D
ClientSampleId :
FB-06272025



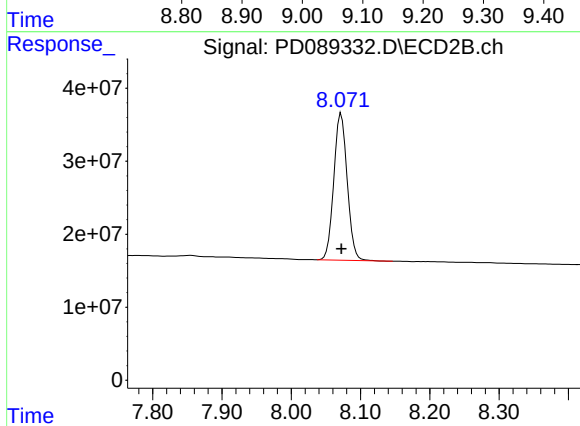
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 356385991
Conc: 21.02 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.075 min
Delta R.T.: 0.003 min
Response: 51727903
Conc: 13.17 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 270856201
Conc: 13.70 ng/ml



CALIBRATION SUMMARY

Lab Name:	<u>Alliance</u>	Contract:	<u>CAMP02</u>	
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q2458</u>	
Instrument ID:	<u>ECD_D</u>	Calibration Date(s):	<u>06/17/2025</u>	<u>06/17/2025</u>
		Calibration Times:	<u>15:52</u>	<u>16:47</u>

LAB FILE ID:	RT 100 =	<u>PD088993.D</u>	RT 075 =	<u>PD088994.D</u>
RT 050 = PD088995.D	RT 025 =	PD088996.D	RT 005 =	PD088997.D

[illegible]

Calibration Times: 15:52 16:47

LAB FILE ID:	RT 100 =	<u>PD088993.D</u>	RT 075 =	<u>PD088994.D</u>
	RT 050 =	PD088995.D	RT 025 =	PD088996.D
			RT 005 =	PD088997.D

[illegible]



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CALIBRATION FACTOR OF INITIAL CALIBRATION

Lab Name: Alliance
Lab Code: ACE
Instrument ID: ECD_D

Contract: CAMP02
SDG NO.: Q2458

Calibration Date(s): 06/17/2025 06/17/2025
Calibration Times: 15:52 16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

LAB FILE ID:				CF 100 =	<u>PD088993.D</u>	CF 075 =	<u>PD088994.D</u>
CF 050 =		<u>PD088995.D</u>	CF 025 =		<u>PD088996.D</u>	CF 005 =	<u>PD088997.D</u>
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	3557450000	3439230000	3414120000	3228670000	3520780000	3432050000	4
4,4'-DDE	4564130000	4401240000	4312380000	4092310000	4435110000	4361030000	4
4,4'-DDT	3928430000	3802000000	3763430000	3580670000	3866490000	3788210000	3
Aldrin	5504150000	5312640000	5254800000	4989300000	5504000000	5312980000	4
alpha-BHC	6295370000	6031790000	5874750000	5417910000	5525530000	5829070000	6
alpha-Chlordane	4885060000	4758900000	4734630000	4584390000	5229700000	4838540000	5
beta-BHC	2136750000	2100860000	2137680000	2144670000	2515730000	2207140000	8
Decachlorobiphenyl	3598890000	3628150000	3773200000	3897960000	4739940000	3927630000	12
delta-BHC	5546280000	5292010000	5182350000	4751190000	4892690000	5132900000	6
Dieldrin	4933650000	4797080000	4750270000	4527510000	4978810000	4797470000	4
Endosulfan I	4511200000	4405770000	4412600000	4289200000	4918470000	4507450000	5
Endosulfan II	3829260000	3982180000	4002080000	3816490000	4553000000	4036600000	7
Endosulfan sulfate	3786900000	3718160000	3741600000	3667440000	4271990000	3837220000	6
Endrin	4238140000	4109540000	4096430000	3878120000	4305700000	4125590000	4
Endrin aldehyde	2964260000	2935670000	2964230000	2966830000	3488720000	3063940000	8
Endrin ketone	4050100000	3970810000	4004900000	3895960000	4443590000	4073070000	5
gamma-BHC (Lindane)	5900830000	5686510000	5566590000	5221690000	5532800000	5581680000	4
gamma-Chlordane	4869330000	4801300000	4716110000	4508810000	5045020000	4788110000	4
Heptachlor	5621840000	5429540000	5358570000	5116310000	5733550000	5451960000	4
Heptachlor epoxide	4794850000	4673610000	4749650000	4554690000	5346870000	4823940000	6
Methoxychlor	1926110000	1918890000	1971480000	1982080000	2304580000	2020630000	8
Tetrachloro-m-xylene	2812300000	2762040000	2767280000	2751820000	3172230000	2853130000	6



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CALIBRATION FACTOR OF INITIAL CALIBRATION

Lab Name: Alliance
Lab Code: ACE
Instrument ID: ECD_D

Contract: CAMP02
SDG NO.: Q2458

Calibration Date(s): 06/17/2025 06/17/2025
Calibration Times: 15:52 16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

LAB FILE ID:		CF 100 =	<u>PD088993.D</u>	CF 075 =	<u>PD088994.D</u>		
CF 050 =		<u>PD088995.D</u>	CF 025 =	<u>PD088996.D</u>	CF 005 =	<u>PD088997.D</u>	
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	17754100000	17899700000	18191100000	18805800000	22996200000	19129400000	11
4,4'-DDE	21343500000	21440200000	21823600000	22564100000	27409000000	22916100000	11
4,4'-DDT	19391000000	19327800000	19594100000	20005200000	23078500000	20279300000	8
Aldrin	22719200000	22747700000	23159400000	23852800000	28607200000	24217300000	10
alpha-BHC	25649000000	25438700000	25849200000	26153100000	30895500000	26797100000	9
alpha-Chlordane	21191300000	21145300000	21570900000	22517200000	28059300000	22896800000	13
beta-BHC	9903630000	9918550000	10198200000	10624100000	13086800000	10746200000	12
Decachlorobiphenyl	17946500000	17939700000	18428600000	19427500000	25107500000	19770000000	15
delta-BHC	23614400000	23533700000	23918700000	24298300000	29003000000	24873600000	9
Dieldrin	21438700000	21622000000	22120600000	22941800000	27919200000	23208500000	12
Endosulfan I	19170300000	19201000000	19315500000	20752900000	25772600000	20842500000	14
Endosulfan II	18473200000	18573800000	19087100000	20002300000	24846800000	20196600000	13
Endosulfan sulfate	17925600000	18043500000	18492500000	19409200000	24234500000	19621100000	13
Endrin	20535400000	19906600000	20644700000	21276500000	26206500000	21713900000	12
Endrin aldehyde	13773600000	13920200000	14379400000	15195500000	19287900000	15311300000	15
Endrin ketone	19587400000	19824900000	20446500000	21535300000	26614600000	21601700000	13
gamma-BHC (Lindane)	23622800000	23442000000	23749900000	24221500000	28748400000	24756900000	9
gamma-Chlordane	22229100000	22022500000	22519500000	23493700000	29348800000	23922700000	13
Heptachlor	23291600000	23365200000	23922200000	24716000000	29896800000	25038400000	11
Heptachlor epoxide	20080600000	20251900000	20827300000	21695300000	26635200000	21898100000	12
Methoxychlor	9712450000	9897880000	10259100000	10830900000	13140100000	10768100000	13
Tetrachloro-m-xylene	15879800000	15866400000	15978000000	16730900000	20310700000	16953200000	11



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INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Contract: CAMP02

Lab Code: CHEM Case No.: Q2458 SAS No.: Q2458 SDG NO.: Q2458

Instrument ID: ECD_D Date(s) Analyzed: 06/17/2025 06/17/2025

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	AMOUNT (ng)	PEAK	RT	RT WINDOW		CALIBRATION FACTOR
				FROM	TO	
Toxaphene	500	1	6.24	6.14	6.34	33601800
		2	6.44	6.34	6.54	47528600
		3	7.15	7.05	7.25	87292700
		4	7.56	7.46	7.66	111357000
		5	7.93	7.83	8.03	63112600



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INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Contract: CAMP02

Lab Code: CHEM Case No.: Q2458 SAS No.: Q2458 SDG NO.: Q2458

Instrument ID: ECD_D Date(s) Analyzed: 06/17/2025 06/17/2025

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	AMOUNT (ng)	PEAK	RT	RT WINDOW		CALIBRATION FACTOR
				FROM	TO	
Toxaphene	500	1	5.47	5.37	5.57	158160000
		2	5.65	5.55	5.75	108820000
		3	6.76	6.66	6.86	511856000
		4	7.20	7.10	7.30	350145000
		5	7.33	7.23	7.43	253920000

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088993.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 15:52
 Operator : AR\AJ
 Sample : PSTDICC100
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDICC100

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:22:36 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:21:59 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	281.2E6	1588.0E6	101.627	99.386
28)	SA Decachlor...	9.072	8.072	359.9E6	1794.6E6	95.380	97.384
Target Compounds							
2)	A alpha-BHC	3.999	3.393	629.5E6	2564.9E6	107.160	99.226
3)	MA gamma-BHC...	4.330	3.729	590.1E6	2362.3E6	106.005	99.465
4)	MA Heptachlor	4.929	4.083	562.2E6	2329.2E6	104.913	97.364
5)	MB Aldrin	5.271	4.369	550.4E6	2271.9E6	104.745	98.099
6)	B beta-BHC	4.515	4.025	213.7E6	990.4E6	99.957	97.111
7)	B delta-BHC	4.764	4.262	554.6E6	2361.4E6	107.022	98.728
8)	B Heptachlo...	5.691	4.873	479.5E6	2008.1E6	100.952	96.415
9)	A Endosulfan I	6.074	5.247	451.1E6	1917.0E6	102.235	99.248
10)	B gamma-Chl...	5.946	5.126	486.9E6	2222.9E6	103.249	98.710
11)	B alpha-Chl...	6.027	5.191	488.5E6	2119.1E6	103.177	98.241
12)	B 4,4'-DDE	6.195	5.375	456.4E6	2134.3E6	105.838	97.800
13)	MA Dieldrin	6.347	5.513	493.4E6	2143.9E6	103.861	96.917
14)	MA Endrin	6.574	5.789	423.8E6	2053.5E6	103.459	99.471
15)	B Endosulfa...	6.786	6.081	382.9E6	1847.3E6	95.682	96.784
16)	A 4,4'-DDD	6.705	5.930	355.7E6	1775.4E6	104.198	97.597
17)	MA 4,4'-DDT	7.020	6.184	392.8E6	1939.1E6	104.384	98.963
18)	B Endrin al...	6.915	6.259	296.4E6	1377.4E6	100.001	95.787
19)	B Endosulfa...	7.148	6.482	378.7E6	1792.6E6	101.211	96.935
20)	A Methoxychlor	7.493	6.755	192.6E6	971.2E6	97.699	94.672
21)	B Endrin ke...	7.629	6.992	405.0E6	1958.7E6	101.129	95.798
22)	Mirex	8.113	7.186	280.1E6	1495.3E6	96.467	95.966

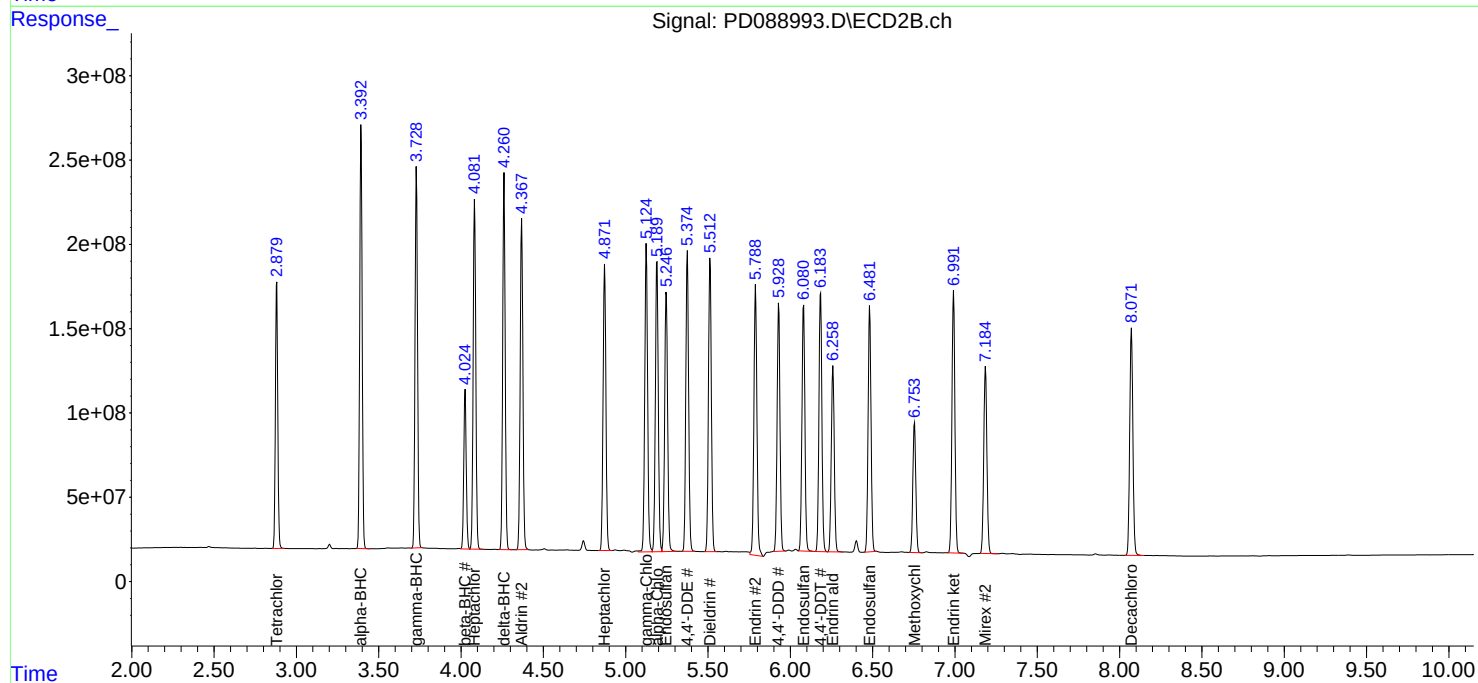
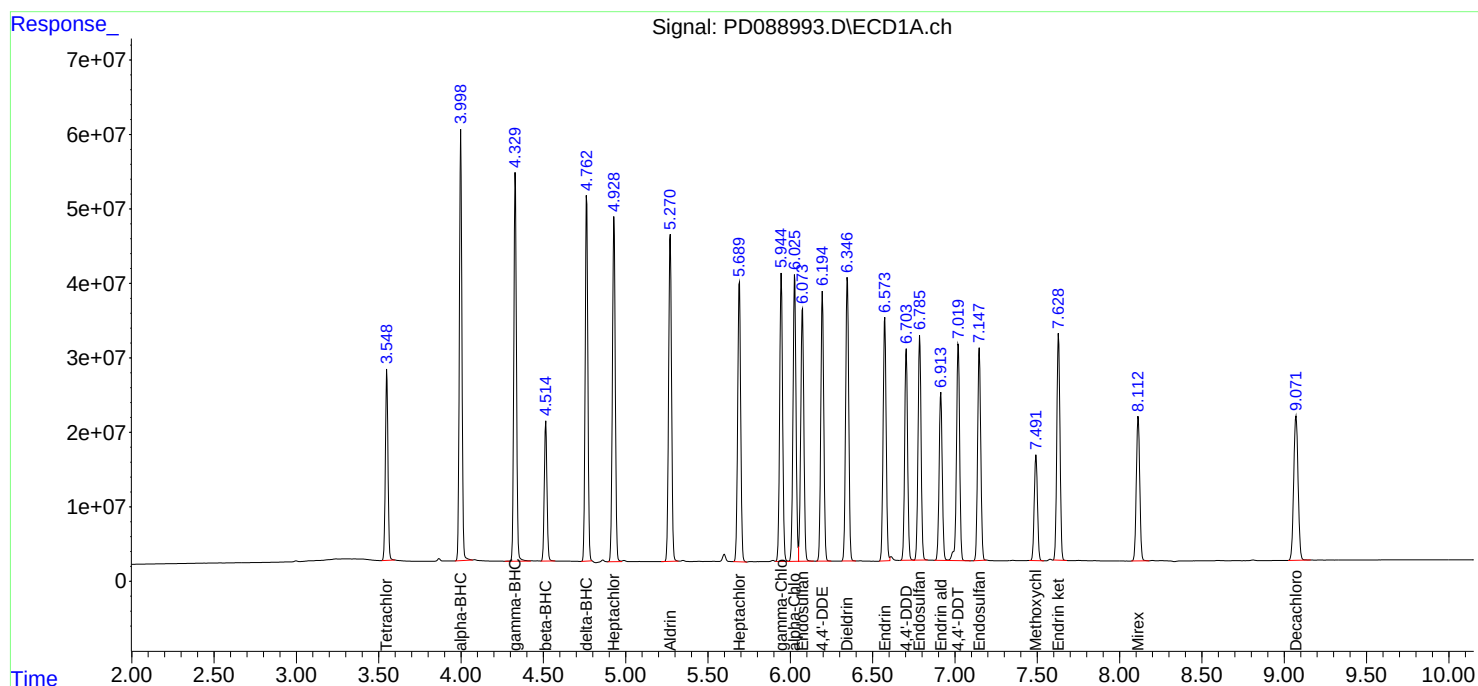
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088993.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 15:52
Operator : AR\AJ
Sample : PSTDICC100
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PSTDICC100

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:22:36 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:21:59 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088994.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 16:06
 Operator : AR\AJ
 Sample : PSTDICC075
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDICC075

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:22:55 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:21:59 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.880	207.2E6	1190.0E6	74.858	74.476
2)	SA Decachlor...	9.072	8.072	272.1E6	1345.5E6	72.117	73.010
Target Compounds							
2)	A alpha-BHC	3.999	3.393	452.4E6	1907.9E6	77.005	73.809
3)	MA gamma-BHC...	4.330	3.729	426.5E6	1758.1E6	76.616	74.028
4)	MA Heptachlor	4.929	4.082	407.2E6	1752.4E6	75.993	73.254
5)	MB Aldrin	5.271	4.369	398.4E6	1706.1E6	75.826	73.667
6)	B beta-BHC	4.515	4.025	157.6E6	743.9E6	73.708	72.943
7)	B delta-BHC	4.764	4.262	396.9E6	1765.0E6	76.587	73.793
8)	B Heptachlo...	5.690	4.872	350.5E6	1518.9E6	73.799	72.928
9)	A Endosulfan I	6.074	5.247	330.4E6	1440.1E6	74.884	74.556
10)	B gamma-Chl...	5.945	5.125	360.1E6	1651.7E6	76.355	73.345
11)	B alpha-Chl...	6.026	5.190	356.9E6	1585.9E6	75.384	73.520
12)	B 4,4'-DDE	6.195	5.375	330.1E6	1608.0E6	76.545	73.682
13)	MA Dieldrin	6.346	5.513	359.8E6	1621.7E6	75.739	73.309
14)	MA Endrin	6.574	5.789	308.2E6	1493.0E6	75.240	72.319
15)	B Endosulfa...	6.785	6.080	298.7E6	1393.0E6	74.627	72.983
16)	A 4,4'-DDD	6.704	5.929	257.9E6	1342.5E6	75.552	73.799
17)	MA 4,4'-DDT	7.020	6.183	285.1E6	1449.6E6	75.769	73.981
18)	B Endrin al...	6.914	6.259	220.2E6	1044.0E6	74.277	72.605
19)	B Endosulfa...	7.149	6.482	278.9E6	1353.3E6	74.530	73.179
20)	A Methoxychlor	7.493	6.754	143.9E6	742.3E6	73.000	72.360
21)	B Endrin ke...	7.629	6.991	297.8E6	1486.9E6	74.361	72.720
22)	Mirex	8.113	7.185	210.7E6	1130.6E6	72.582	72.559

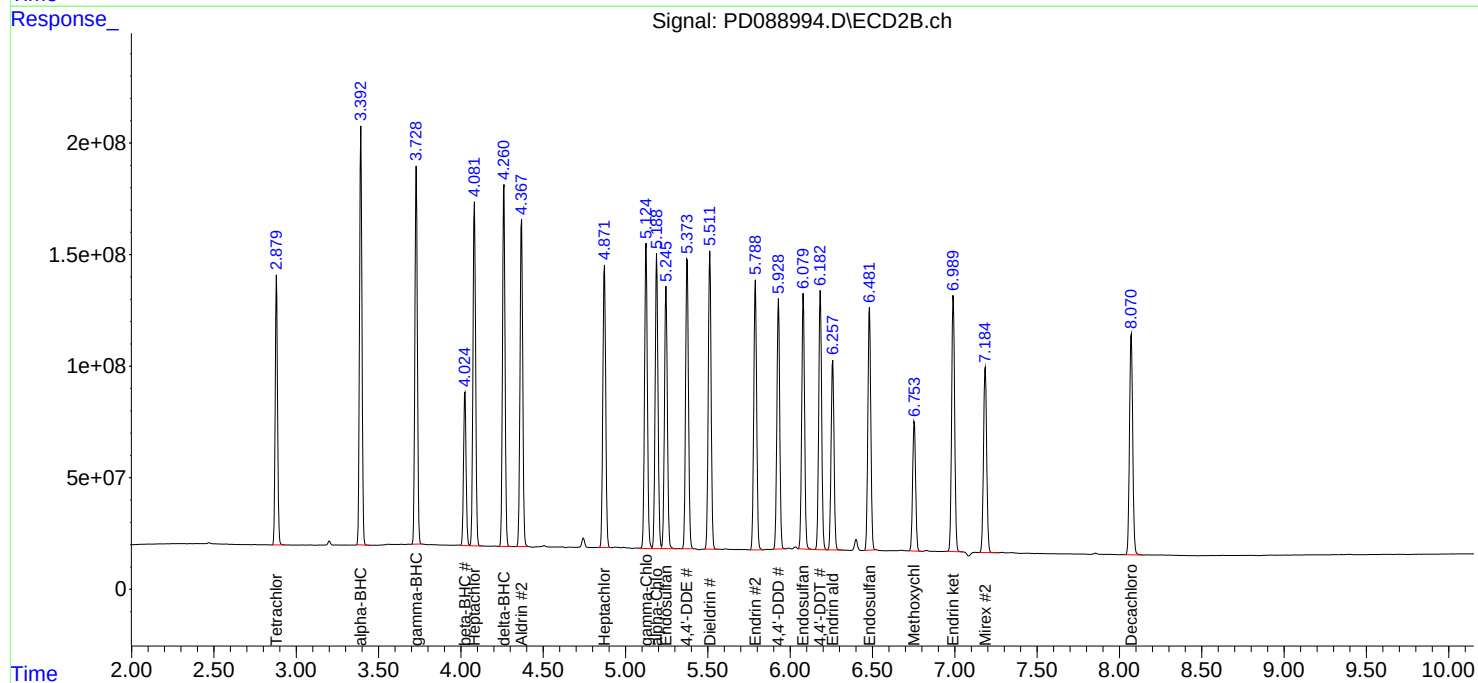
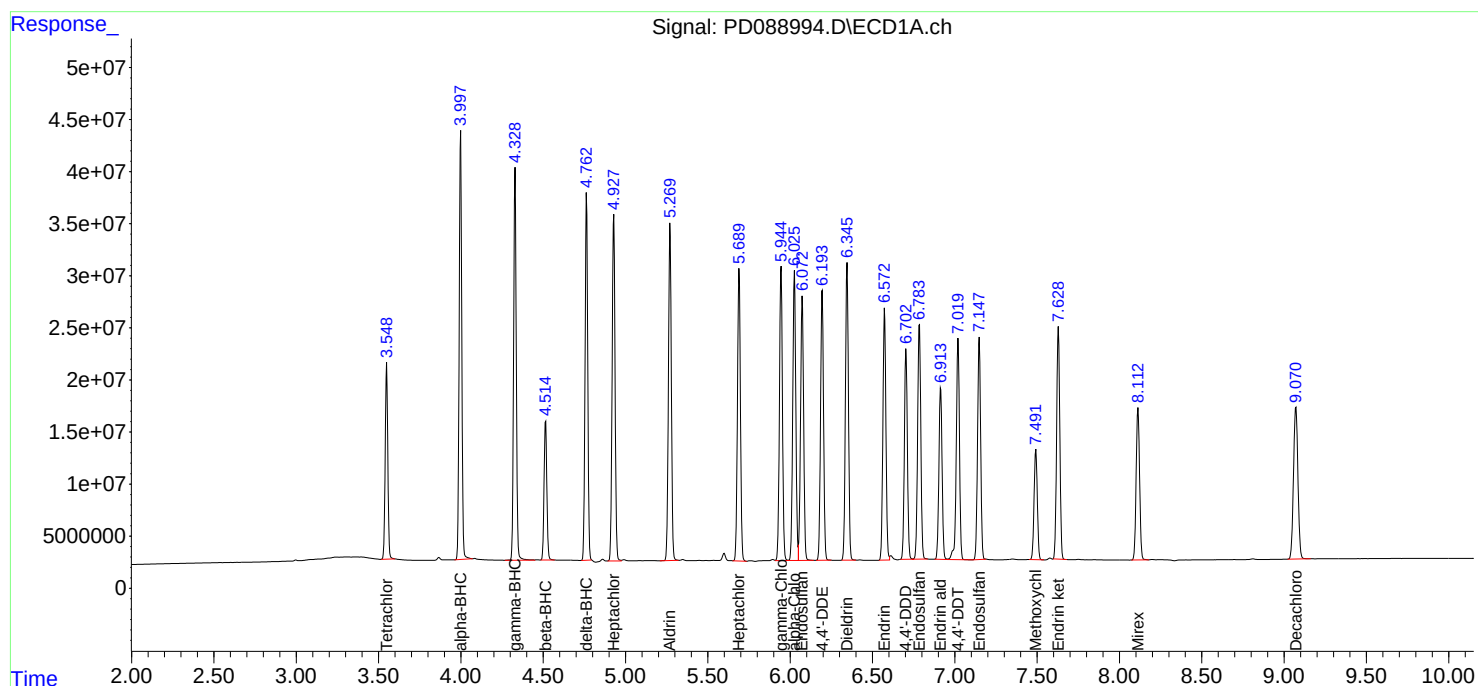
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088994.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 16:06
Operator : AR\AJ
Sample : PSTDICC075
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PSTDICC075

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:22:55 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:21:59 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088995.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 16:20
 Operator : AR\AJ
 Sample : PSTDICC050
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDICC050

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:23:15 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:21:59 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.880	138.4E6	798.9E6	50.000	50.000
28)	SA Decachlor...	9.072	8.072	188.7E6	921.4E6	50.000	50.000
Target Compounds							
2)	A alpha-BHC	3.999	3.393	293.7E6	1292.5E6	50.000	50.000
3)	MA gamma-BHC...	4.330	3.730	278.3E6	1187.5E6	50.000	50.000
4)	MA Heptachlor	4.929	4.083	267.9E6	1196.1E6	50.000	50.000
5)	MB Aldrin	5.271	4.369	262.7E6	1158.0E6	50.000	50.000
6)	B beta-BHC	4.515	4.026	106.9E6	509.9E6	50.000	50.000
7)	B delta-BHC	4.764	4.262	259.1E6	1195.9E6	50.000	50.000
8)	B Heptachlo...	5.691	4.873	237.5E6	1041.4E6	50.000	50.000
9)	A Endosulfan I	6.075	5.247	220.6E6	965.8E6	50.000	50.000
10)	B gamma-Chl...	5.946	5.126	235.8E6	1126.0E6	50.000	50.000
11)	B alpha-Chl...	6.027	5.191	236.7E6	1078.5E6	50.000	50.000
12)	B 4,4'-DDE	6.196	5.375	215.6E6	1091.2E6	50.000	50.000
13)	MA Dieldrin	6.347	5.513	237.5E6	1106.0E6	50.000	50.000
14)	MA Endrin	6.575	5.789	204.8E6	1032.2E6	50.000	50.000
15)	B Endosulfa...	6.786	6.081	200.1E6	954.4E6	50.000	50.000
16)	A 4,4'-DDD	6.705	5.930	170.7E6	909.6E6	50.000	50.000
17)	MA 4,4'-DDT	7.021	6.184	188.2E6	979.7E6	50.000	50.000
18)	B Endrin al...	6.915	6.259	148.2E6	719.0E6	50.000	50.000
19)	B Endosulfa...	7.149	6.483	187.1E6	924.6E6	50.000	50.000
20)	A Methoxychlor	7.493	6.755	98574001	513.0E6	50.000	50.000
21)	B Endrin ke...	7.630	6.992	200.2E6	1022.3E6	50.000	50.000
22)	Mirex	8.114	7.186	145.2E6	779.1E6	50.000	50.000

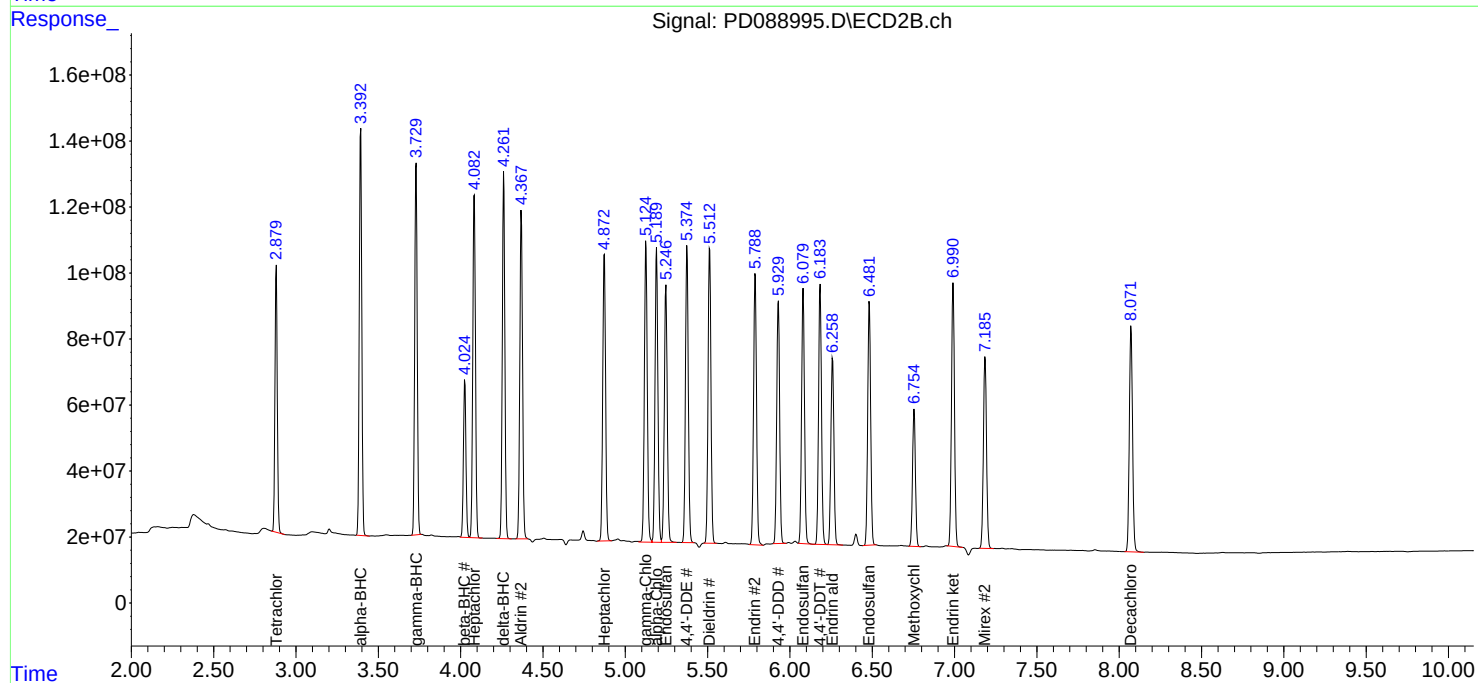
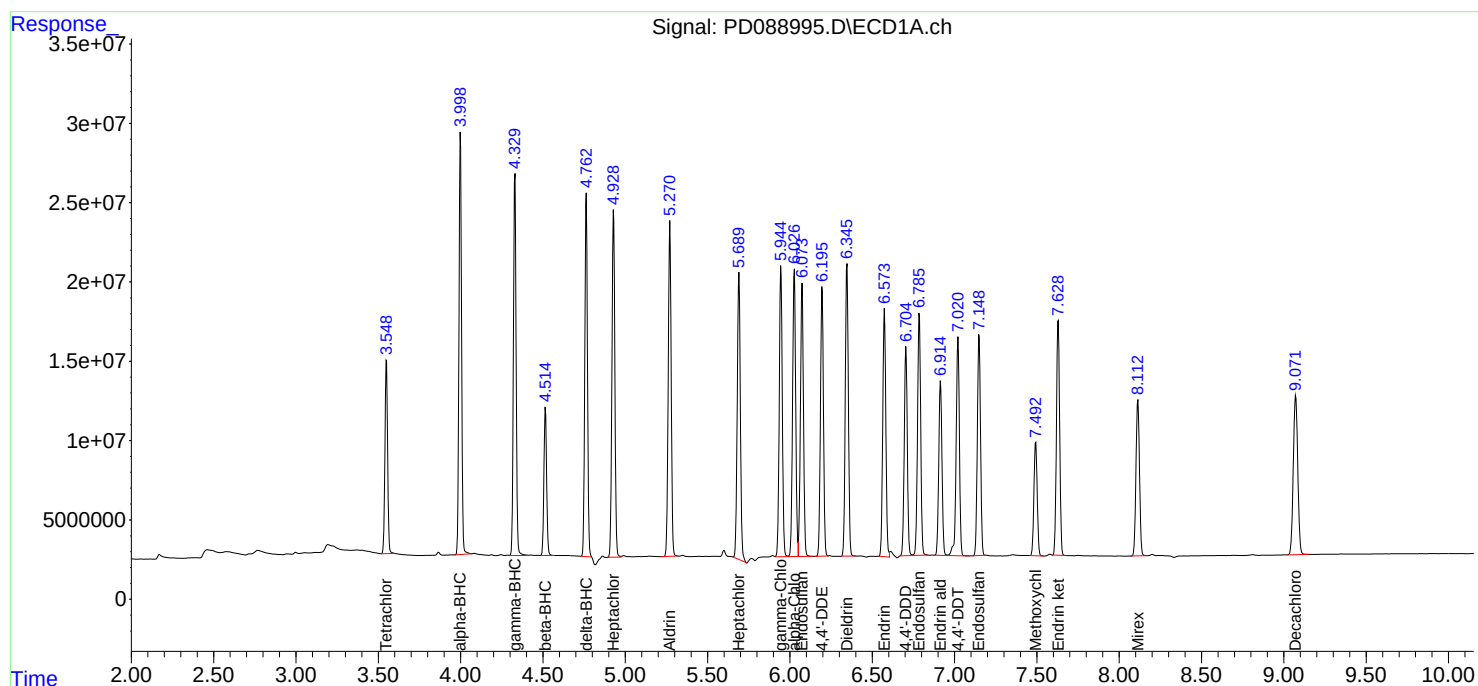
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088995.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 16:20
Operator : AR\AJ
Sample : PSTDICC050
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PSTDICC050

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:23:15 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:21:59 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088996.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 16:33
 Operator : AR\AJ
 Sample : PSTDICC025
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDICC025

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 06/18/2025
 Supervised By :mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:23:35 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:21:59 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.881	68795424	418.3E6	24.860	26.178
28)	SA Decachlor...	9.072	8.072	97448897	485.7E6	25.827	26.355
Target Compounds							
2)	A alpha-BHC	3.999	3.393	135.4E6	653.8E6	23.056	25.294
3)	MA gamma-BHC...	4.330	3.729	130.5E6	605.5E6	23.451	25.496
4)	MA Heptachlor	4.929	4.083	127.9E6	617.9E6	23.870	25.830
5)	MB Aldrin	5.271	4.368	124.7E6	596.3E6	23.737	25.748
6)	B beta-BHC	4.515	4.025	53616740	265.6E6	25.082	26.044
7)	B delta-BHC	4.763	4.261	118.8E6	607.5E6	22.920	25.397
8)	B Heptachlo...	5.690	4.872	113.9E6	542.4E6	23.974	26.042
9)	A Endosulfan I	6.074	5.247	107.2E6	518.8E6	24.301	26.860
10)	B gamma-Chl...	5.945	5.125	112.7E6	587.3E6	23.901	26.082
11)	B alpha-Chl...	6.026	5.190	114.6E6	562.9E6	24.207	26.097
12)	B 4,4'-DDE	6.195	5.375	102.3E6	564.1E6	23.724	25.848
13)	MA Dieldrin	6.346	5.513	113.2E6	573.5E6	23.828	25.928
14)	MA Endrin	6.574	5.787	96953033	531.9E6	23.668	25.765m
15)	B Endosulfa...	6.786	6.080	95412213	500.1E6	23.841	26.199
16)	A 4,4'-DDD	6.704	5.929	80716755	470.1E6	23.642	25.845
17)	MA 4,4'-DDT	7.020	6.183	89516720	500.1E6	23.786	25.525
18)	B Endrin al...	6.914	6.258	74170726	379.9E6	25.022	26.419
19)	B Endosulfa...	7.148	6.481	91685883	485.2E6	24.504	26.239
20)	A Methoxychlor	7.492	6.753	49551931	270.8E6	25.134	26.394
21)	B Endrin ke...	7.629	6.990	97399015	538.4E6	24.320	26.331
22)	Mirex	8.113	7.184	74764205	414.2E6	25.753	26.583

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088996.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 16:33
Operator : AR\AJ
Sample : PSTDICC025
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDICC025

Manual Integrations

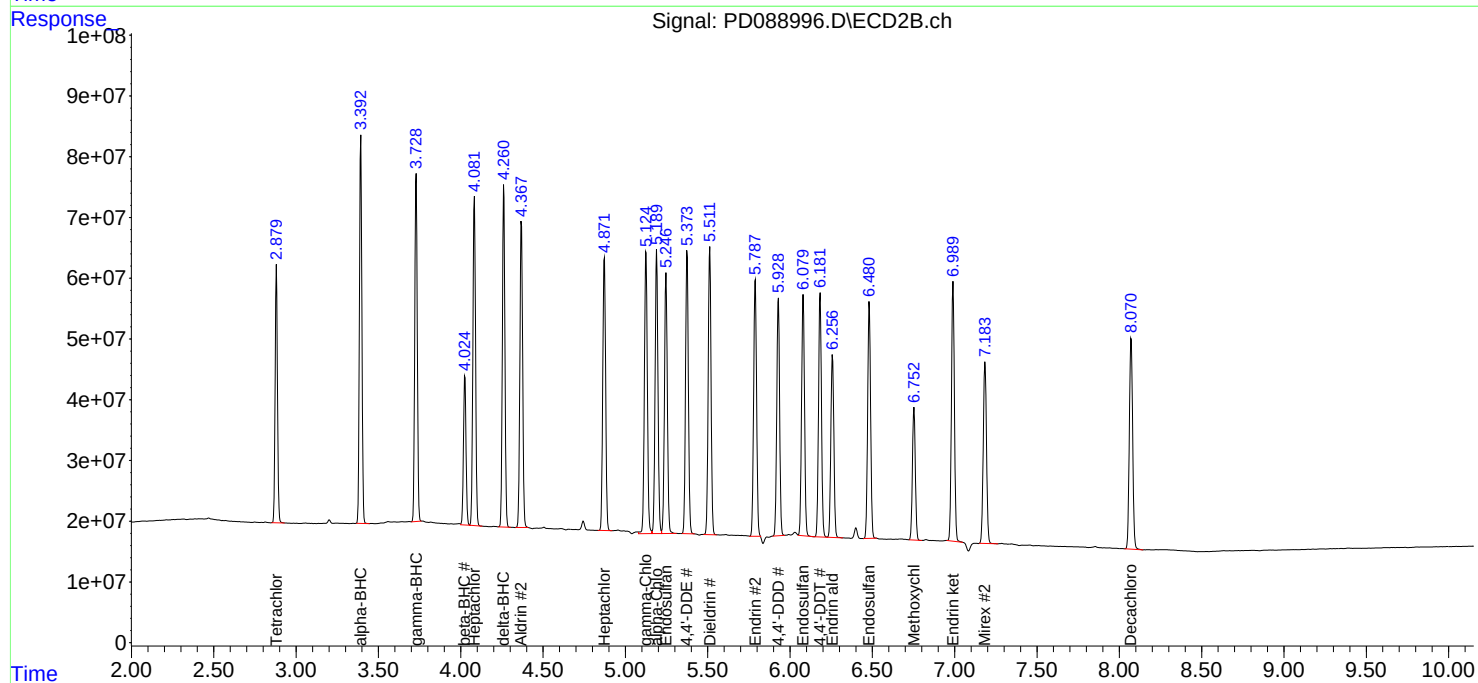
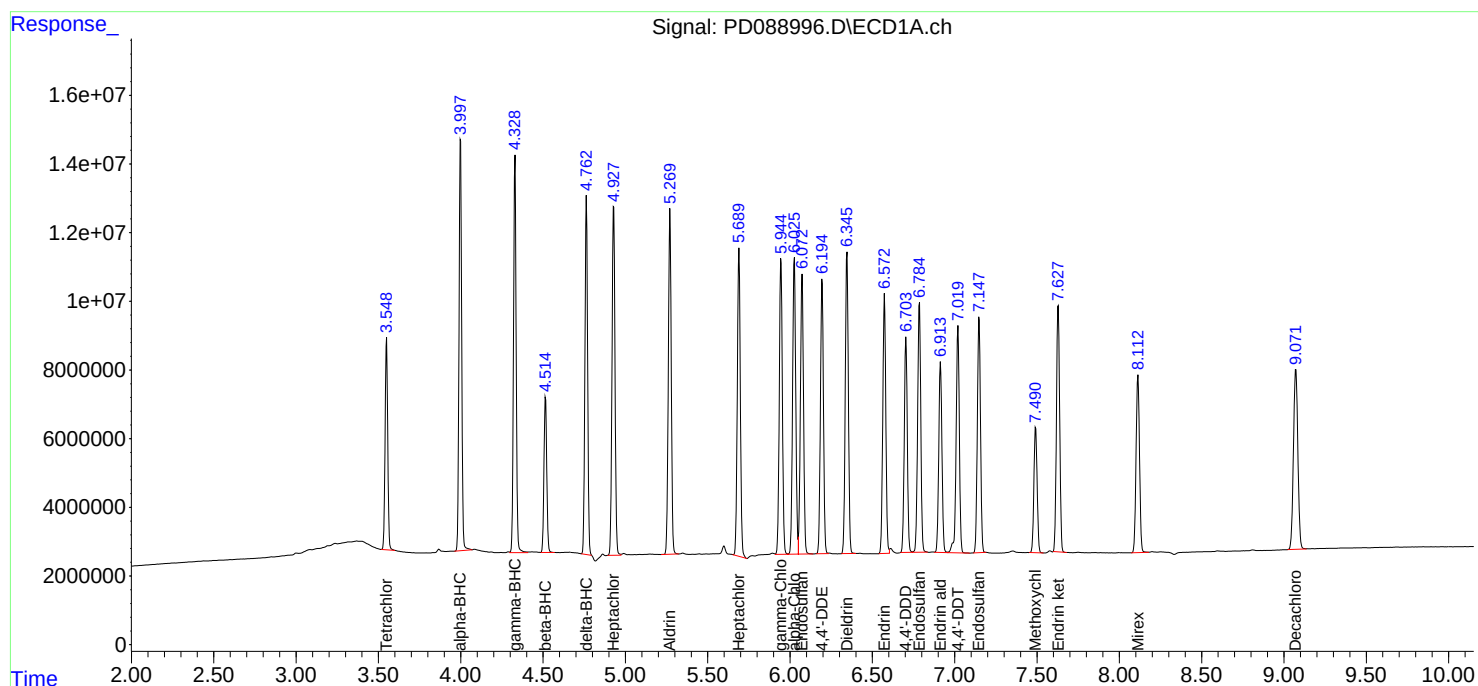
APPROVED

Reviewed By :Abdul Mirza 06/18/2025

Supervised By :mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:23:35 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:21:59 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088997.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 16:47
 Operator : AR\AJ
 Sample : PSTDICC005
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 06/18/2025
 Supervised By :mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:23:55 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:21:59 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.549	2.880	15861136	101.6E6	5.732	6.356
2) SA Decachlor...	9.073	8.072	23699682	125.5E6	6.281	6.812
Target Compounds						
2) A alpha-BHC	3.999	3.393	27627643	154.5E6	4.703	5.976 #
3) MA gamma-BHC...	4.330	3.729	27664017	143.7E6	4.970	6.052
4) MA Heptachlor	4.929	4.083	28667725	149.5E6	5.350	6.249
5) MB Aldrin	5.270	4.368	27520008	143.0E6	5.237	6.176
6) B beta-BHC	4.515	4.025	12578663	65433774	5.884	6.416
7) B delta-BHC	4.763	4.262	24463430	145.0E6	4.721	6.063 #
8) B Heptachlo...	5.690	4.873	26734361	133.2E6	5.629	6.394
9) A Endosulfan I	6.074	5.247	24592339	128.9E6	5.573	6.671
10) B gamma-Chl...	5.946	5.126	25225085	146.7E6	5.349	6.516
11) B alpha-Chl...	6.026	5.190	26148512	140.3E6	5.523	6.504
12) B 4,4'-DDE	6.194	5.375	22175565	137.0E6	5.142	6.280
13) MA Dieldrin	6.346	5.512	24894050	139.6E6	5.241	6.311
14) MA Endrin	6.573	5.787	21528490	131.0E6	5.255	6.347m
15) B Endosulfa...	6.785	6.080	22765018	124.2E6	5.688	6.509
16) A 4,4'-DDD	6.704	5.929	17603896	115.0E6	5.156	6.321
17) MA 4,4'-DDT	7.020	6.183	19332472	115.4E6	5.137	5.889
18) B Endrin al...	6.914	6.259	17443585	96439397	5.885	6.707
19) B Endosulfa...	7.149	6.482	21359951	121.2E6	5.709	6.553
20) A Methoxychlor	7.493	6.754	11522878	65700483	5.845	6.404
21) B Endrin ke...	7.630	6.990	22217957	133.1E6	5.548	6.508m
22) Mirex	8.113	7.185	18484356	106.1E6	6.367	6.807

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088997.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 16:47
Operator : AR\AJ
Sample : PSTDICC005
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDICC005

Manual Integrations

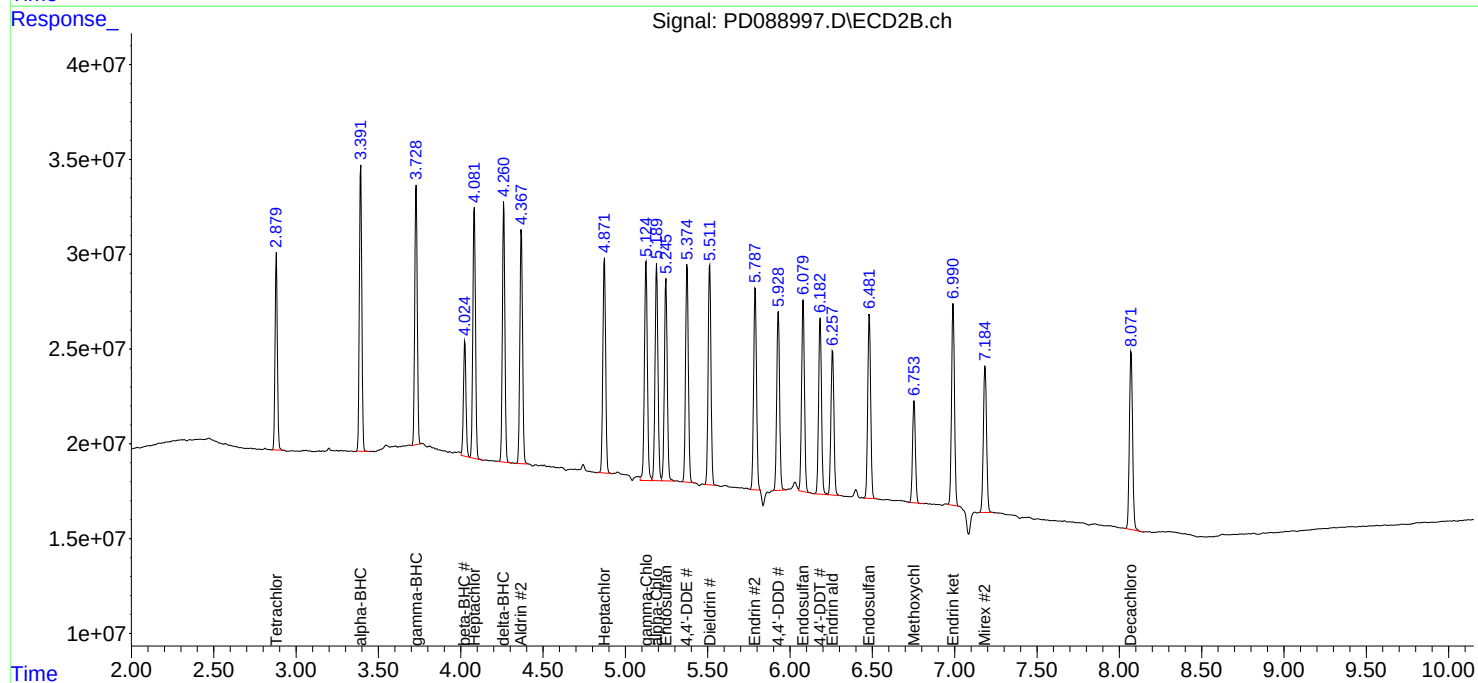
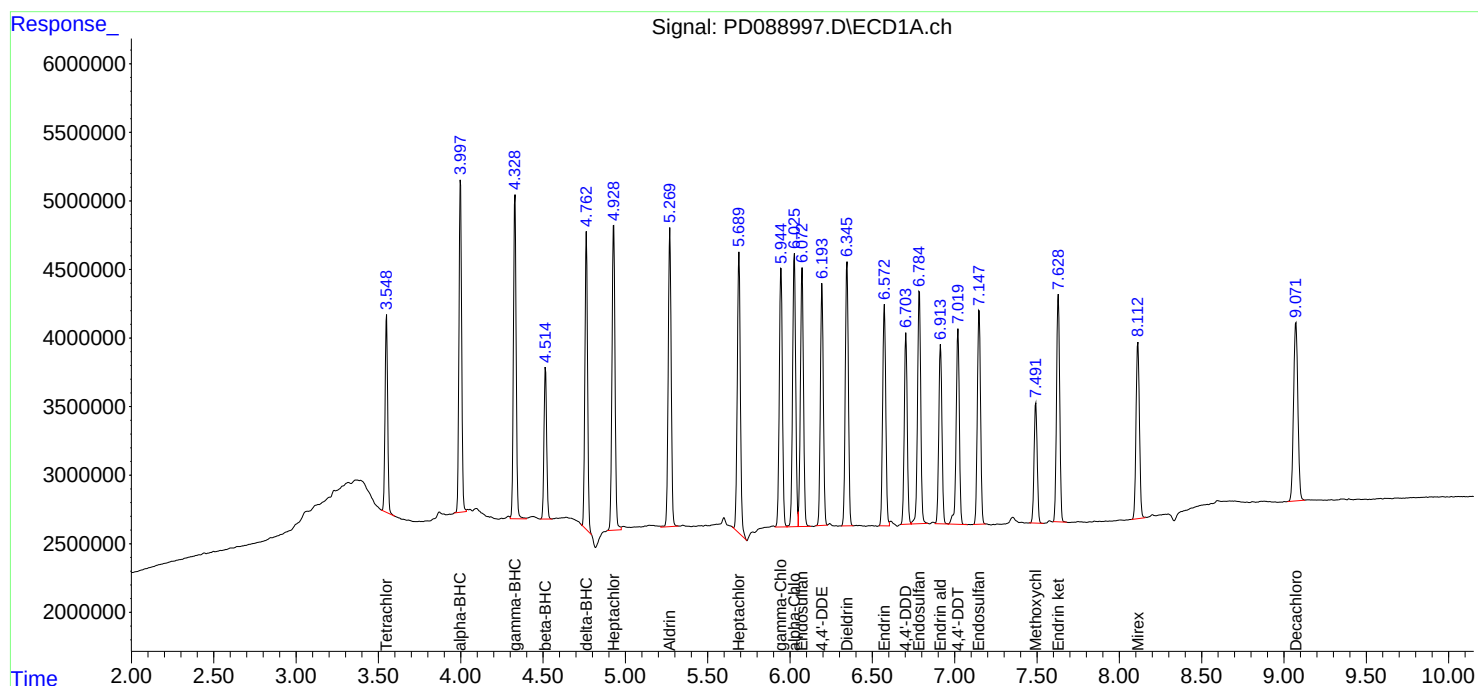
APPROVED

Reviewed By :Abdul Mirza 06/18/2025

Supervised By :mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:23:55 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:21:59 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD089000.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 17:28
 Operator : AR\AJ
 Sample : PCHLORICC500
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PCHLORICC500

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 04:59:08 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 04:57:57 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.880	127.0E6	1003.3E6	50.000	50.000
28)	SA Decachlor...	9.072	8.072	175.3E6	940.3E6	50.000	50.000
Target Compounds							
23)	Chlordane-1	4.715	3.905	103.3E6	448.7E6	500.000	500.000
24)	Chlordane-2	5.241	4.487	103.7E6	451.9E6	500.000	500.000
25)	Chlordane-3	5.946	5.126	425.8E6	1413.2E6	500.000	500.000
26)	Chlordane-4	6.032	5.190	511.2E6	1191.6E6	500.000	500.000
27)	Chlordane-5	6.870	6.090	88549137	546.3E6	500.000	500.000

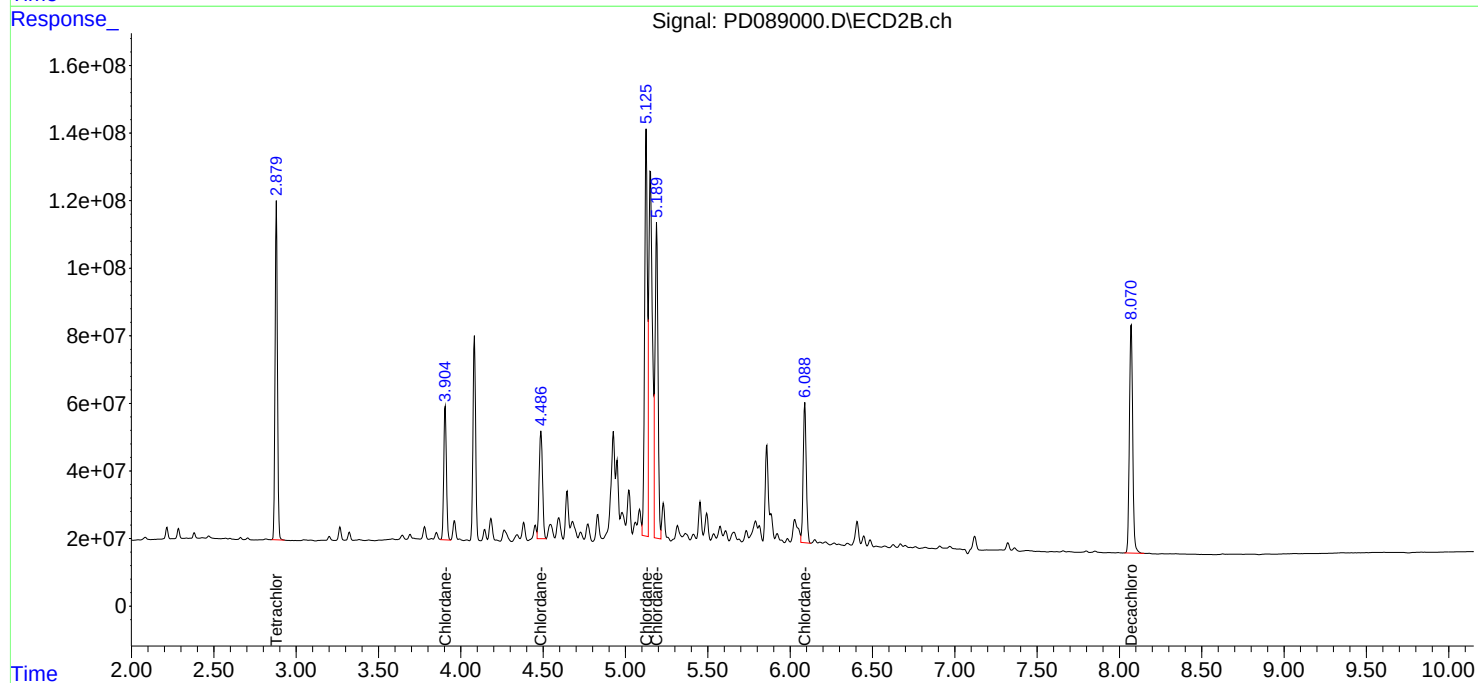
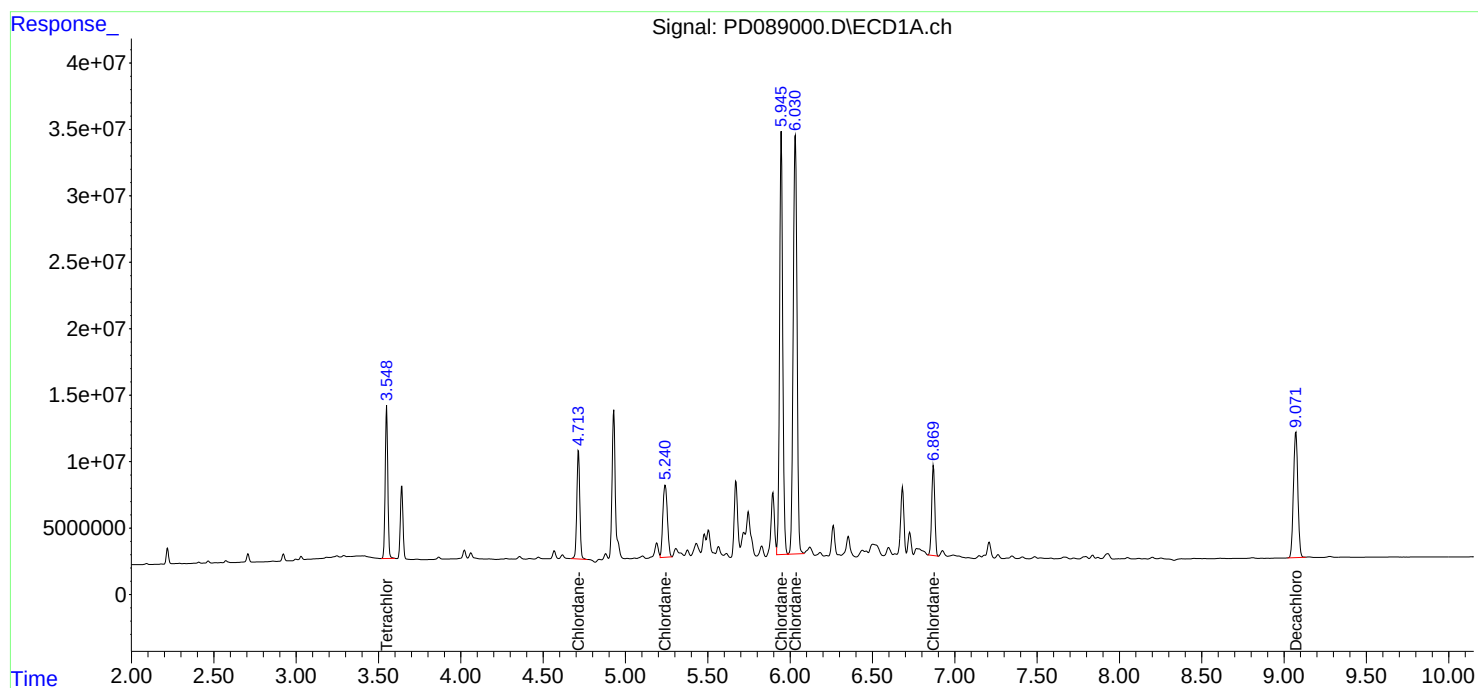
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD089000.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 17:28
Operator : AR\AJ
Sample : PCHLORICC500
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PCHLORICC500

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 04:59:08 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 04:57:57 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD089005.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 18:36
 Operator : AR\AJ
 Sample : PTOXICC500
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PTOXICC500

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 04:30:48 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\DTX061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 04:29:28 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x 0. Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

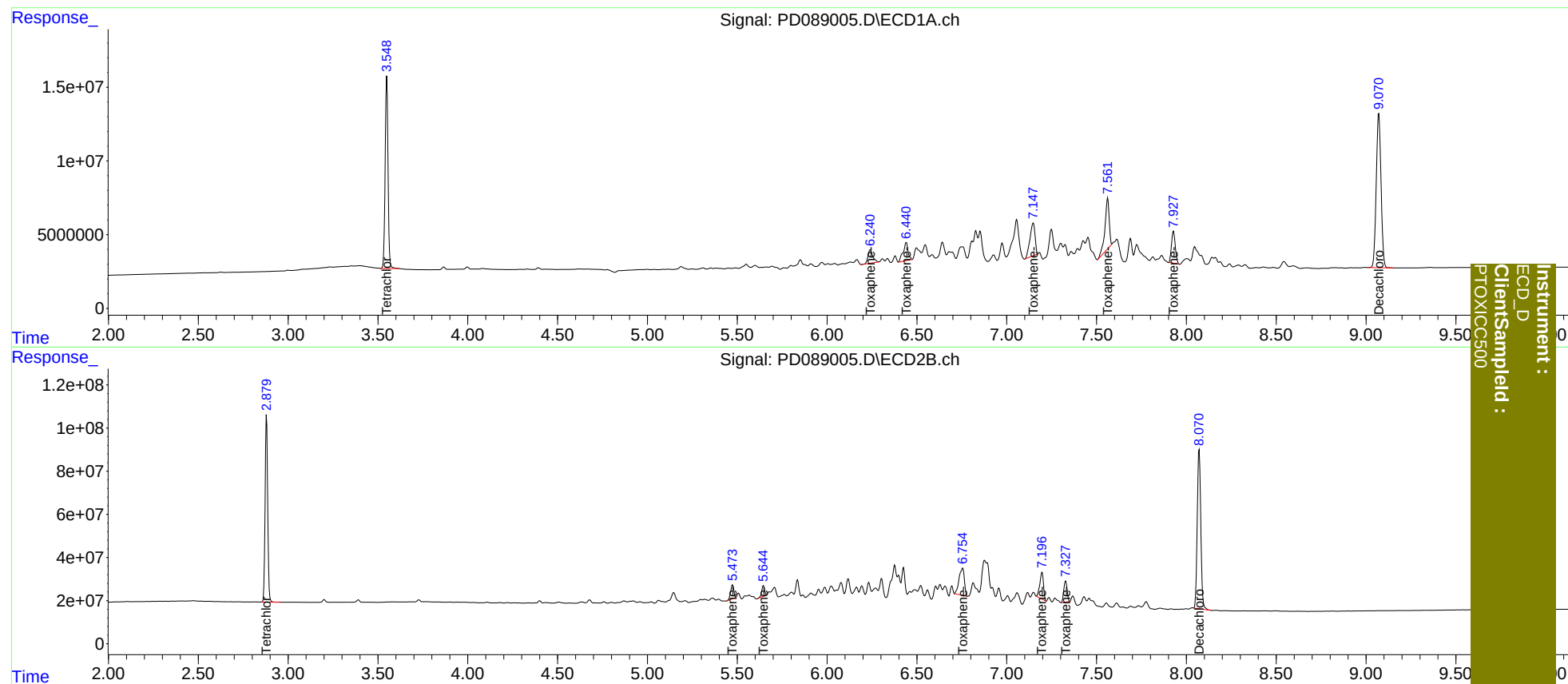
System Monitoring Compounds						
1) SA Tetrachlo...	3.549	2.880	143.6E6	868.2E6	50.000	50.000
7) SA Decachlor...	9.071	8.071	194.1E6	1009.6E6	50.000	50.000
Target Compounds						
2) Toxaphene-1	6.241	5.474	16800895	79079940	500.000	500.000
3) Toxaphene-2	6.441	5.646	23764296	54409973	500.000	500.000
4) Toxaphene-3	7.148	6.755	43646368	255.9E6	500.000	500.000
5) Toxaphene-4	7.562	7.197	55678343	175.1E6	500.000	500.000
6) Toxaphene-5	7.928	7.328	31556310	127.0E6	500.000	500.000

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD089005.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 18:36
Operator : AR\AJ
Sample : PTOXICC500
Misc :
ALS Vial : 17 Sample Multiplier: 1

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 04:30:48 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\DTX061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 04:29:28 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0. Signal #2 Info : 30M x 0.32mm x0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD089008.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 19:17
 Operator : AR\AJ
 Sample : PSTDICV050
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 ICPD061825

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:46:15 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.880	139.5E6	848.9E6	48.909	50.072
28)	SA Decachlor...	9.072	8.073	188.5E6	996.5E6	47.986	50.403
Target Compounds							
2)	A alpha-BHC	3.999	3.393	295.8E6	1350.4E6	50.737	50.394
3)	MA gamma-BHC...	4.330	3.729	280.5E6	1249.3E6	50.254	50.462
4)	MA Heptachlor	4.929	4.082	270.9E6	1252.2E6	49.682	50.010
5)	MB Aldrin	5.271	4.368	265.0E6	1218.5E6	49.876	50.317
6)	B beta-BHC	4.515	4.025	107.7E6	537.8E6	48.800	50.043
7)	B delta-BHC	4.764	4.262	261.5E6	1252.0E6	50.946	50.333
8)	B Heptachlo...	5.691	4.872	235.4E6	1093.7E6	48.806	49.946
9)	A Endosulfan I	6.075	5.247	222.0E6	1045.8E6	49.252	50.175
10)	B gamma-Chl...	5.946	5.126	237.3E6	1184.2E6	49.570	49.500
11)	B alpha-Chl...	6.027	5.190	238.3E6	1135.5E6	49.257	49.590
12)	B 4,4'-DDE	6.195	5.375	217.6E6	1145.1E6	49.901	49.970
13)	MA Dieldrin	6.347	5.513	238.9E6	1159.0E6	49.793	49.940
14)	MA Endrin	6.574	5.790	203.1E6	1100.0E6	49.223	50.657
15)	B Endosulfa...	6.786	6.081	195.6E6	1002.6E6	48.453	49.643
16)	A 4,4'-DDD	6.705	5.930	170.0E6	955.9E6	49.526	49.972
17)	MA 4,4'-DDT	7.021	6.184	188.0E6	1027.1E6	49.627	50.650
18)	B Endrin al...	6.915	6.259	149.8E6	759.8E6	48.894	49.621
19)	B Endosulfa...	7.149	6.483	187.3E6	976.2E6	48.824	49.755
20)	A Methoxychlor	7.493	6.755	98271542	536.8E6	48.634	49.853
21)	B Endrin ke...	7.630	6.992	200.8E6	1076.8E6	49.300	49.848
22)	Mirex	8.114	7.186	147.3E6	825.2E6	48.452	49.481

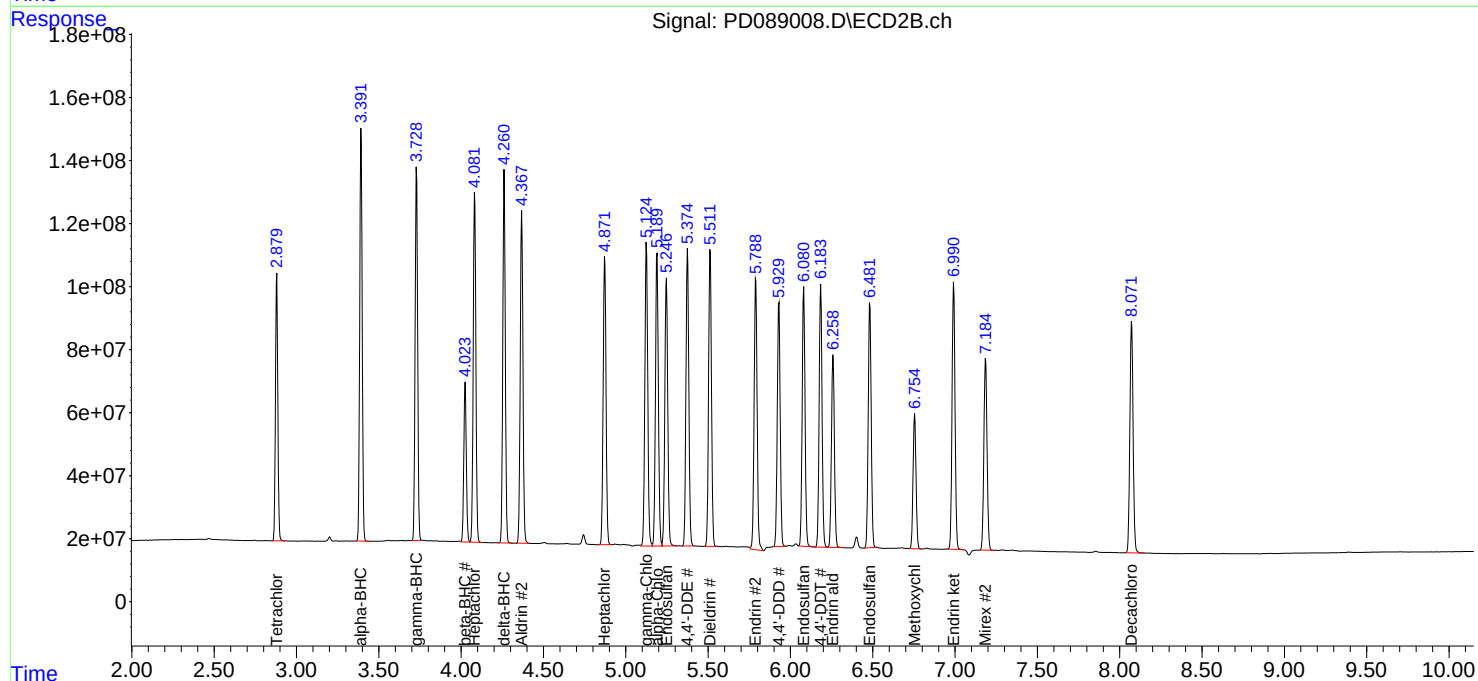
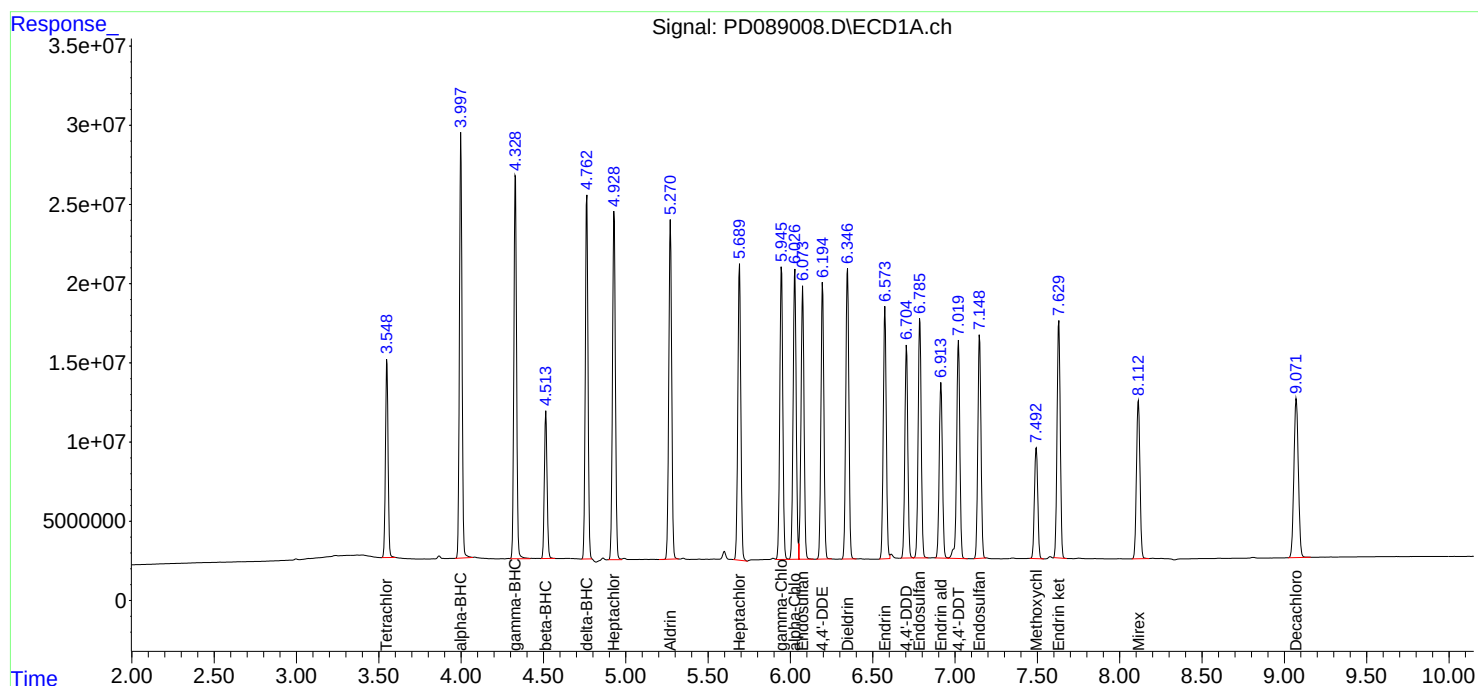
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD089008.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 19:17
Operator : AR\AJ
Sample : PSTDICV050
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
ICVPD061825

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:46:15 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD089009.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 19:31
 Operator : AR\AJ
 Sample : PCHLORICV500
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 ICVPD061825CHLOR

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:06:39 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:05:26 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.880	127.5E6	1002.8E6	48.435	48.165
28)	SA Decachlor...	9.072	8.072	176.1E6	945.9E6	47.752	47.396
Target Compounds							
23)	Chlordane-1	4.714	3.905	104.1E6	451.3E6	486.162	484.057
24)	Chlordane-2	5.241	4.487	104.8E6	451.6E6	480.042	474.511
25)	Chlordane-3	5.946	5.125	422.4E6	1412.7E6	489.742	475.621
26)	Chlordane-4	6.031	5.189	512.9E6	1195.8E6	488.156	477.027
27)	Chlordane-5	6.870	6.089	88790524	548.7E6	485.833	484.867

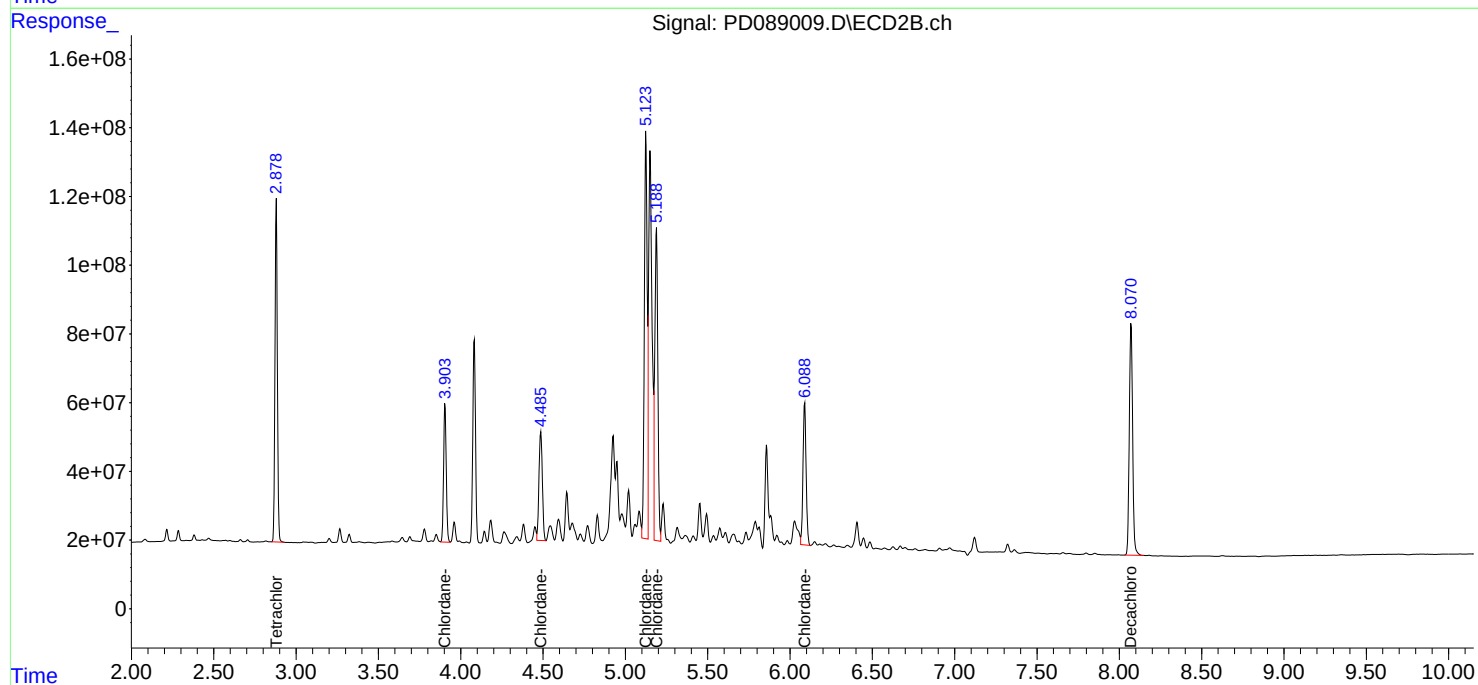
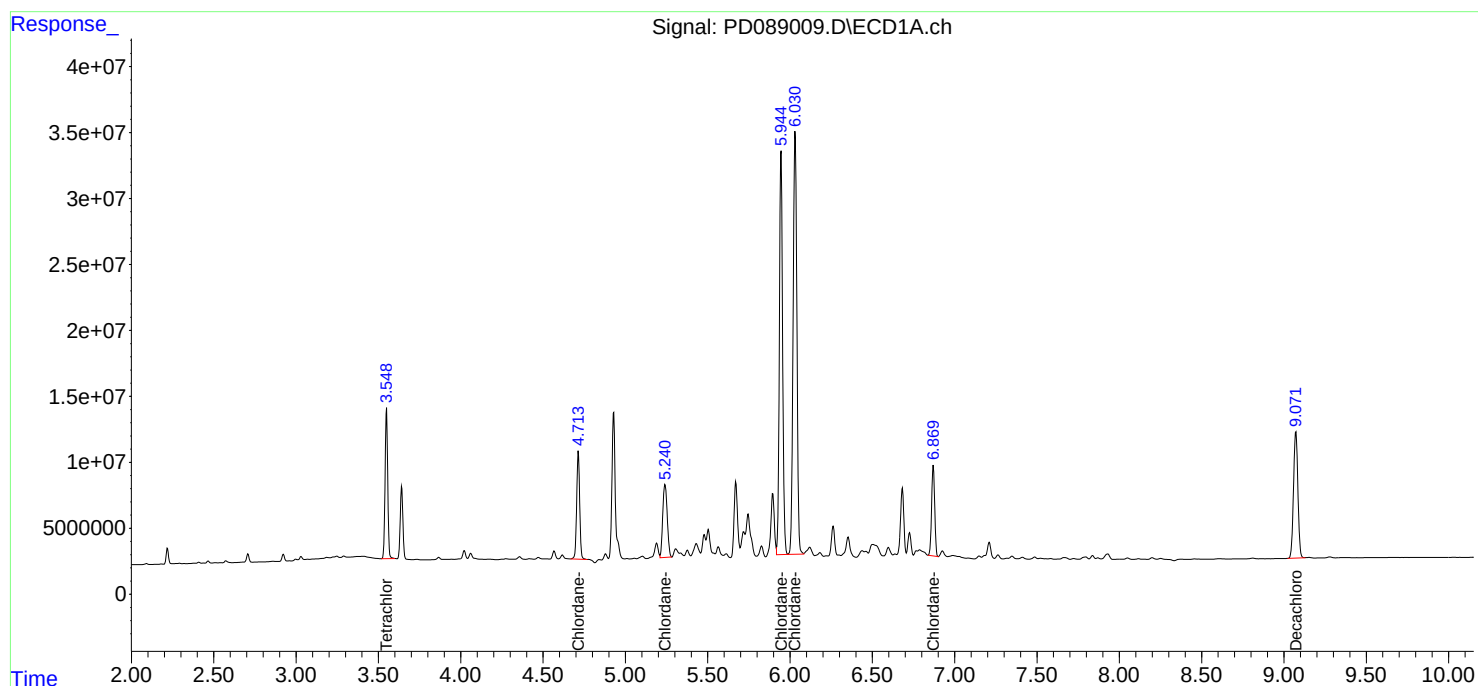
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD089009.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 19:31
Operator : AR\AJ
Sample : PCHLORICV500
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
ICVPD061825CHLOR

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:06:39 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:05:26 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD089010.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 19:44
 Operator : AR\AJ
 Sample : PTOXICV500
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 ICVPD061825TOX

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 04:37:01 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\DTX061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 04:35:42 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x 0. Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

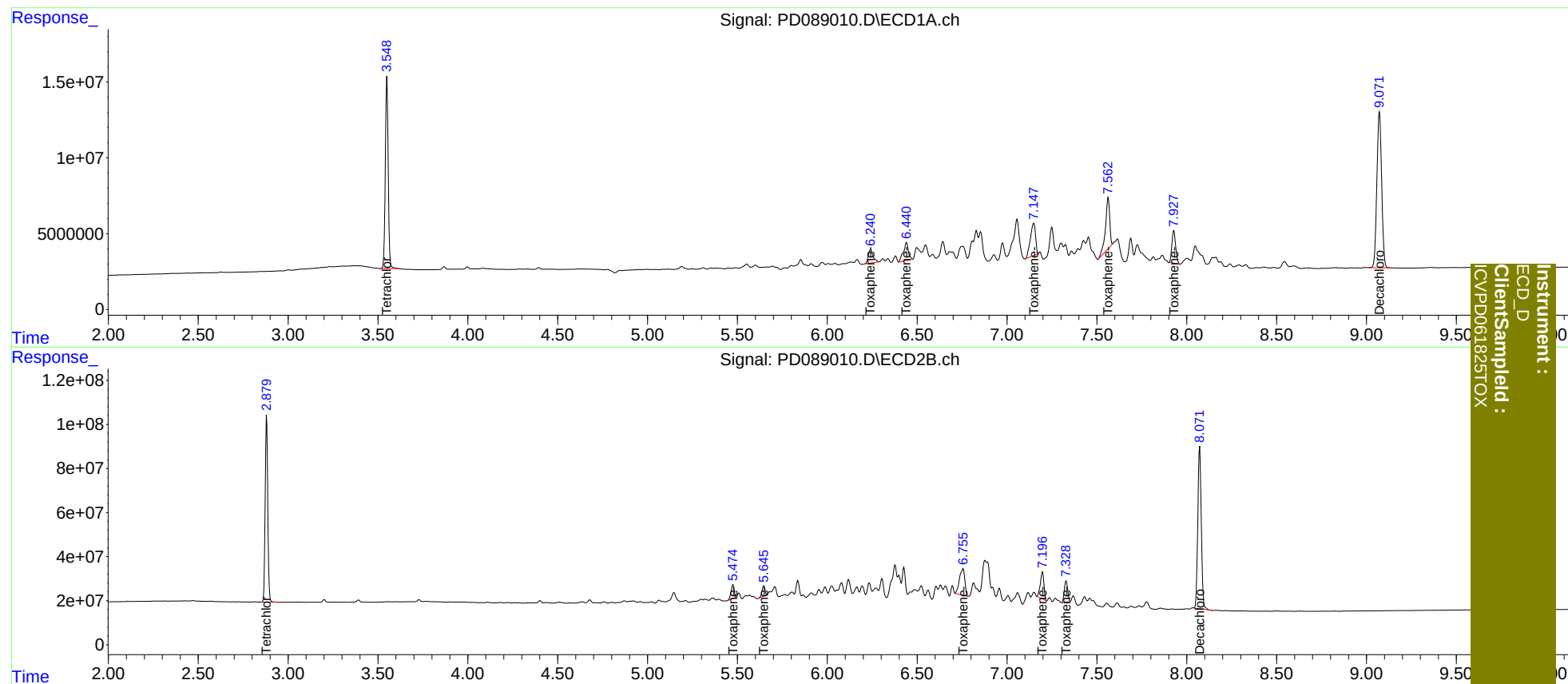
System Monitoring Compounds						
1) SA Tetrachlo...	3.550	2.881	140.5E6	849.8E6	49.126	48.618
7) SA Decachlor...	9.072	8.072	190.0E6	991.7E6	48.552	48.009
Target Compounds						
2) Toxaphene-1	6.241	5.475	16556024	76482668	496.105	468.012
3) Toxaphene-2	6.441	5.647	23088671	52728004	503.218	488.483
4) Toxaphene-3	7.149	6.756	42578482	247.9E6	493.980	497.712
5) Toxaphene-4	7.563	7.198	54552672	171.5E6	492.551	493.627
6) Toxaphene-5	7.928	7.329	31161300	123.5E6	501.825	508.354

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD089010.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 19:44
Operator : AR\AJ
Sample : PTOXICV500
Misc :
ALS Vial : 22 Sample Multiplier: 1

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 04:37:01 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\DTX061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 04:35:42 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0. Signal #2 Info : 30M x 0.32mm x0.25µm





284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 13:35

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.08	9.07	8.97	9.17	-0.01
Tetrachloro-m-xylene	3.56	3.55	3.45	3.65	-0.01
alpha-BHC	4.01	4.00	3.90	4.10	-0.01
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.77	4.76	4.66	4.86	-0.01
gamma-BHC (Lindane)	4.34	4.33	4.23	4.43	-0.01
Heptachlor	4.94	4.93	4.83	5.03	-0.01
Aldrin	5.28	5.27	5.17	5.37	-0.01
Heptachlor epoxide	5.70	5.69	5.59	5.79	-0.01
Endosulfan I	6.08	6.08	5.98	6.18	0.00
Dieldrin	6.36	6.35	6.25	6.45	-0.01
4,4'-DDE	6.21	6.20	6.10	6.30	0.00
Endrin	6.58	6.58	6.48	6.68	0.00
Endosulfan II	6.80	6.79	6.69	6.89	-0.01
4,4'-DDD	6.72	6.71	6.61	6.81	-0.01
Endosulfan sulfate	7.16	7.15	7.05	7.25	-0.01
4,4'-DDT	7.03	7.02	6.92	7.12	-0.01
Methoxychlor	7.50	7.49	7.39	7.59	-0.01
Endrin ketone	7.64	7.63	7.53	7.73	-0.01
Endrin aldehyde	6.92	6.92	6.82	7.02	0.00
alpha-Chlordane	6.04	6.03	5.93	6.13	-0.01
gamma-Chlordane	5.96	5.95	5.85	6.05	0.00



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Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 13:35

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.00
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.38	5.38	5.28	5.48	0.00
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.19	6.18	6.08	6.28	0.00
Methoxychlor	6.76	6.76	6.66	6.86	0.00
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.13	5.13	5.03	5.23	0.00



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL01 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089266.D **Time Analyzed:** 13:35

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.715	6.605	6.805	51.660	50.000	3.3
4,4'-DDE	6.205	6.096	6.296	48.860	50.000	-2.3
4,4'-DDT	7.030	6.921	7.121	43.340	50.000	-13.3
Aldrin	5.280	5.171	5.371	51.140	50.000	2.3
alpha-BHC	4.008	3.899	4.099	52.120	50.000	4.2
alpha-Chlordane	6.036	5.927	6.127	49.710	50.000	-0.6
beta-BHC	4.524	4.415	4.615	48.950	50.000	-2.1
Decachlorobiphenyl	9.084	8.972	9.172	48.780	50.000	-2.4
delta-BHC	4.772	4.664	4.864	53.570	50.000	7.1
Dieldrin	6.356	6.247	6.447	50.140	50.000	0.3
Endosulfan I	6.083	5.975	6.175	49.940	50.000	-0.1
Endosulfan II	6.795	6.686	6.886	48.920	50.000	-2.2
Endosulfan sulfate	7.158	7.049	7.249	48.220	50.000	-3.6
Endrin	6.583	6.475	6.675	46.610	50.000	-6.8
Endrin aldehyde	6.923	6.815	7.015	47.590	50.000	-4.8
Endrin ketone	7.639	7.530	7.730	48.720	50.000	-2.6
gamma-BHC (Lindane)	4.339	4.230	4.430	50.900	50.000	1.8
gamma-Chlordane	5.955	5.846	6.046	50.610	50.000	1.2
Heptachlor	4.938	4.829	5.029	49.650	50.000	-0.7
Heptachlor epoxide	5.700	5.591	5.791	49.160	50.000	-1.7
Methoxychlor	7.502	7.393	7.593	49.560	50.000	-0.9
Tetrachloro-m-xylene	3.558	3.450	3.650	51.440	50.000	2.9



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL01 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089266.D **Time Analyzed:** 13:35

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.931	5.830	6.030	47.120	50.000	-5.8
4,4'-DDE	5.376	5.275	5.475	48.960	50.000	-2.1
4,4'-DDT	6.185	6.084	6.284	42.950	50.000	-14.1
Aldrin	4.370	4.269	4.469	49.490	50.000	-1.0
alpha-BHC	3.393	3.293	3.493	50.180	50.000	0.4
alpha-Chlordane	5.192	5.091	5.291	50.230	50.000	0.5
beta-BHC	4.026	3.926	4.126	47.780	50.000	-4.4
Decachlorobiphenyl	8.074	7.972	8.172	48.360	50.000	-3.3
delta-BHC	4.263	4.162	4.362	48.900	50.000	-2.2
Dieldrin	5.512	5.413	5.613	48.830	50.000	-2.3
Endosulfan I	5.248	5.147	5.347	50.360	50.000	0.7
Endosulfan II	6.082	5.981	6.181	48.700	50.000	-2.6
Endosulfan sulfate	6.484	6.383	6.583	48.060	50.000	-3.9
Endrin	5.791	5.689	5.889	47.960	50.000	-4.1
Endrin aldehyde	6.260	6.159	6.359	47.200	50.000	-5.6
Endrin ketone	6.993	6.892	7.092	48.850	50.000	-2.3
gamma-BHC (Lindane)	3.728	3.630	3.830	49.670	50.000	-0.7
gamma-Chlordane	5.125	5.026	5.226	48.580	50.000	-2.8
Heptachlor	4.083	3.983	4.183	47.120	50.000	-5.8
Heptachlor epoxide	4.872	4.773	4.973	49.620	50.000	-0.8
Methoxychlor	6.756	6.655	6.855	49.370	50.000	-1.3
Tetrachloro-m-xylene	2.880	2.780	2.980	51.270	50.000	2.5

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089266.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 13:35
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:45:03 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.558	2.880	146.8E6	869.2E6	51.443	51.270
28)	SA Decachlor...	9.084	8.074	191.6E6	956.2E6	48.783	48.364
Target Compounds							
2)	A alpha-BHC	4.008	3.393	303.8E6	1344.6E6	52.123	50.178
3)	MA gamma-BHC...	4.339	3.728	284.1E6	1229.7E6	50.904	49.670m
4)	MA Heptachlor	4.938	4.083	270.7E6	1179.9E6	49.655	47.124
5)	MB Aldrin	5.280	4.370	271.7E6	1198.6E6	51.135	49.493
6)	B beta-BHC	4.524	4.026	108.0E6	513.5E6	48.946	47.781
7)	B delta-BHC	4.772	4.263	275.0E6	1216.4E6	53.574m	48.903
8)	B Heptachlo...	5.700	4.872	237.1E6	1086.7E6	49.158	49.625m
9)	A Endosulfan I	6.083	5.248	225.1E6	1049.6E6	49.945	50.359
10)	B gamma-Chl...	5.955	5.125	242.3E6	1162.3E6	50.612	48.584m
11)	B alpha-Chl...	6.036	5.192	240.5E6	1150.2E6	49.710	50.235
12)	B 4,4'-DDE	6.205	5.376	213.1E6	1122.0E6	48.859	48.962
13)	MA Dieldrin	6.356	5.512	240.5E6	1133.4E6	50.137	48.834m
14)	MA Endrin	6.583	5.791	192.3E6	1041.3E6	46.608	47.957
15)	B Endosulfa...	6.795	6.082	197.5E6	983.7E6	48.921	48.704
16)	A 4,4'-DDD	6.715	5.931	177.3E6	901.4E6	51.663	47.119
17)	MA 4,4'-DDT	7.030	6.185	164.2E6	871.1E6	43.336	42.954
18)	B Endrin al...	6.923	6.260	145.8E6	722.7E6	47.590m	47.201
19)	B Endosulfa...	7.158	6.484	185.0E6	943.0E6	48.218	48.062
20)	A Methoxychlor	7.502	6.756	100.1E6	531.7E6	49.562	49.374
21)	B Endrin ke...	7.639	6.993	198.5E6	1055.3E6	48.725	48.854
22)	Mirex	8.123	7.186	145.6E6	790.0E6	47.891	47.365m

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089266.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 13:35
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

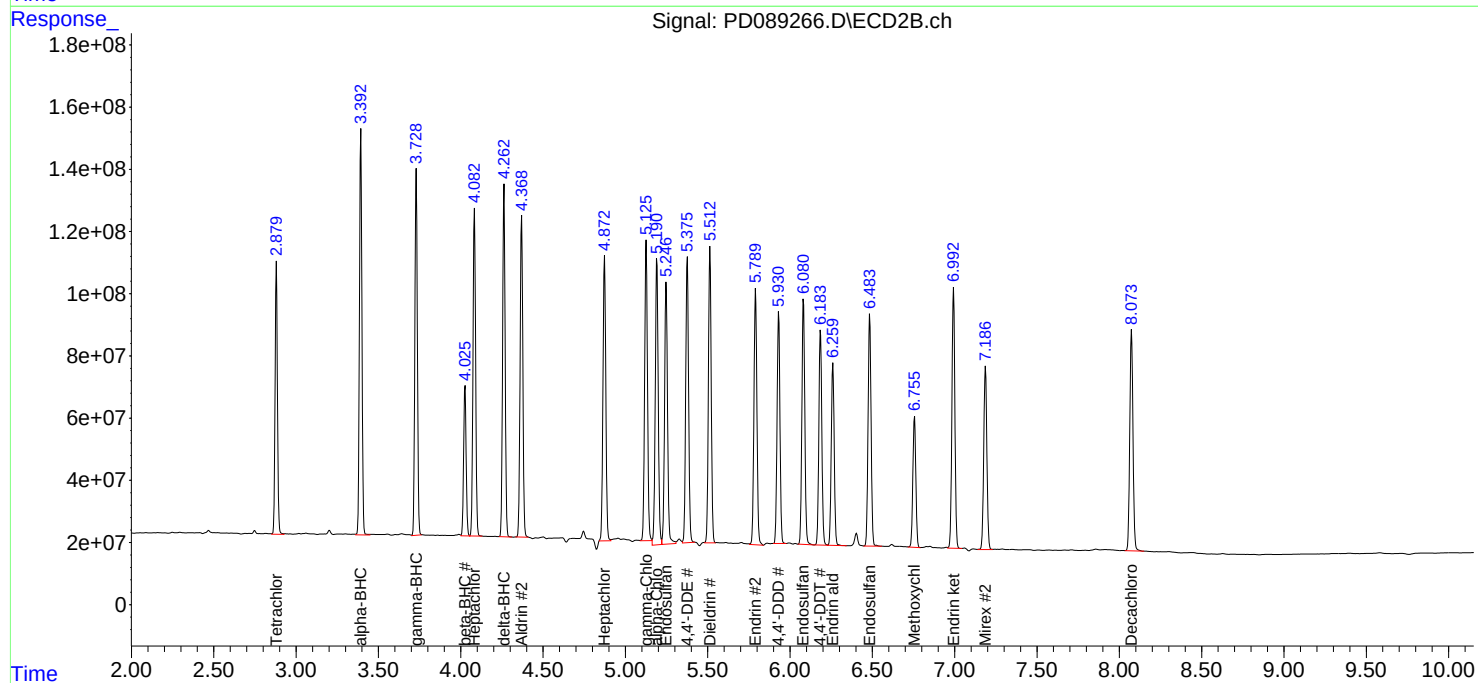
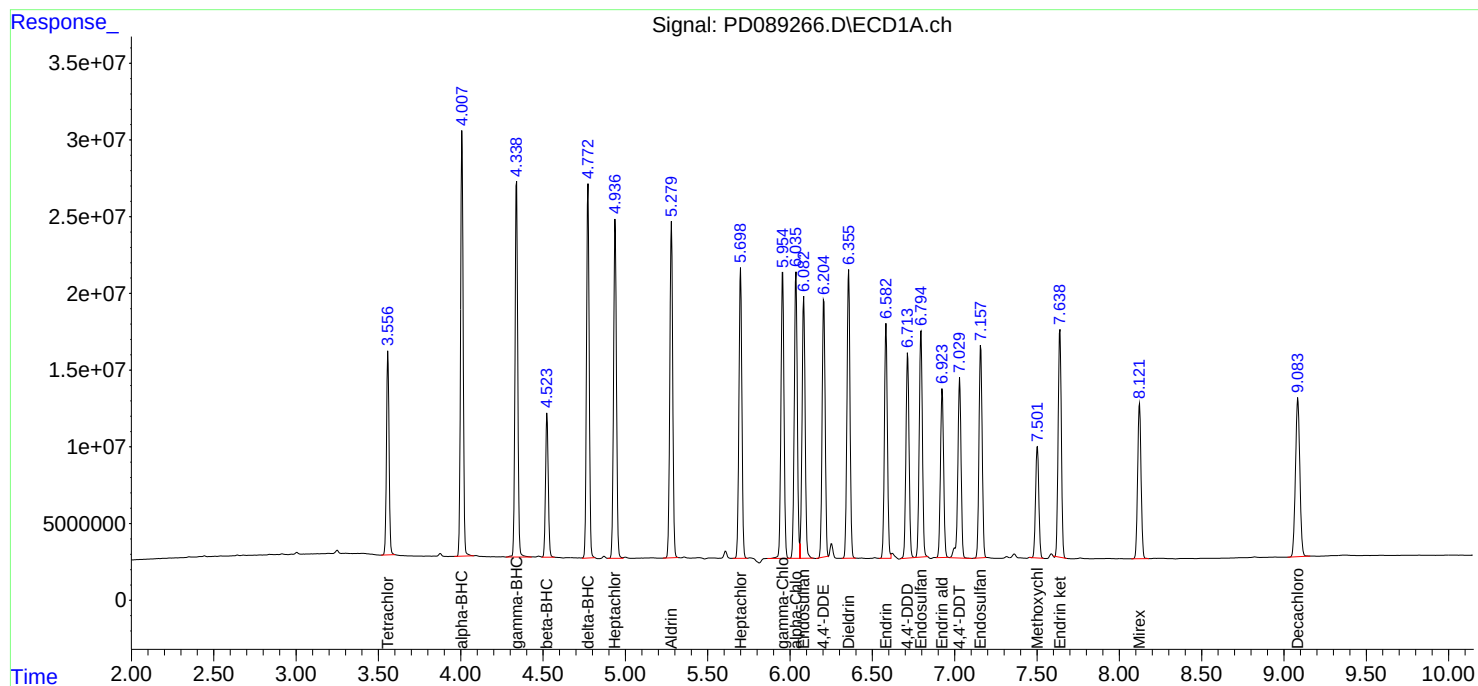
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:45:03 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 16:50

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.07	8.97	9.17	0.00
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	4.00	3.90	4.10	0.00
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.77	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.93	4.83	5.03	0.00
Aldrin	5.27	5.27	5.17	5.37	0.00
Heptachlor epoxide	5.69	5.69	5.59	5.79	0.00
Endosulfan I	6.08	6.08	5.98	6.18	0.01
Dieldrin	6.35	6.35	6.25	6.45	0.00
4,4'-DDE	6.20	6.20	6.10	6.30	0.00
Endrin	6.57	6.58	6.48	6.68	0.01
Endosulfan II	6.79	6.79	6.69	6.89	0.00
4,4'-DDD	6.71	6.71	6.61	6.81	0.00
Endosulfan sulfate	7.15	7.15	7.05	7.25	0.00
4,4'-DDT	7.02	7.02	6.92	7.12	0.00
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.92	6.82	7.02	0.00
alpha-Chlordane	6.03	6.03	5.93	6.13	0.00
gamma-Chlordane	5.95	5.95	5.85	6.05	0.00



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 16:50

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.01
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.38	5.38	5.28	5.48	0.00
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.18	6.18	6.08	6.28	0.00
Methoxychlor	6.75	6.76	6.66	6.86	0.01
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.12	5.13	5.03	5.23	0.01



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group Contract: CAMP02

Lab Code: CHEM SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No.: CCAL02 Date Analyzed: 07/01/2025

Lab Sample No.: PSTDCCC050 Data File : PD089277.D Time Analyzed: 16:50

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.706	6.605	6.805	51.810	50.000	3.6
4,4'-DDE	6.196	6.096	6.296	49.240	50.000	-1.5
4,4'-DDT	7.021	6.921	7.121	44.170	50.000	-11.7
Aldrin	5.272	5.171	5.371	51.690	50.000	3.4
alpha-BHC	4.000	3.899	4.099	52.270	50.000	4.5
alpha-Chlordane	6.028	5.927	6.127	50.240	50.000	0.5
beta-BHC	4.516	4.415	4.615	49.900	50.000	-0.2
Decachlorobiphenyl	9.073	8.972	9.172	48.690	50.000	-2.6
delta-BHC	4.765	4.664	4.864	53.990	50.000	8.0
Dieldrin	6.348	6.247	6.447	50.190	50.000	0.4
Endosulfan I	6.075	5.975	6.175	50.580	50.000	1.2
Endosulfan II	6.787	6.686	6.886	49.450	50.000	-1.1
Endosulfan sulfate	7.150	7.049	7.249	48.310	50.000	-3.4
Endrin	6.574	6.475	6.675	47.220	50.000	-5.6
Endrin aldehyde	6.916	6.815	7.015	48.350	50.000	-3.3
Endrin ketone	7.630	7.530	7.730	49.130	50.000	-1.7
gamma-BHC (Lindane)	4.331	4.230	4.430	51.800	50.000	3.6
gamma-Chlordane	5.947	5.846	6.046	50.960	50.000	1.9
Heptachlor	4.930	4.829	5.029	50.200	50.000	0.4
Heptachlor epoxide	5.691	5.591	5.791	49.950	50.000	-0.1
Methoxychlor	7.493	7.393	7.593	50.180	50.000	0.4
Tetrachloro-m-xylene	3.550	3.450	3.650	51.880	50.000	3.8



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL02 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089277.D **Time Analyzed:** 16:50

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.930	5.830	6.030	48.050	50.000	-3.9
4,4'-DDE	5.375	5.275	5.475	50.250	50.000	0.5
4,4'-DDT	6.183	6.084	6.284	44.120	50.000	-11.8
Aldrin	4.369	4.269	4.469	51.130	50.000	2.3
alpha-BHC	3.394	3.293	3.493	51.500	50.000	3.0
alpha-Chlordane	5.189	5.091	5.291	49.750	50.000	-0.5
beta-BHC	4.026	3.926	4.126	49.490	50.000	-1.0
Decachlorobiphenyl	8.072	7.972	8.172	47.700	50.000	-4.6
delta-BHC	4.263	4.162	4.362	50.640	50.000	1.3
Dieldrin	5.511	5.413	5.613	50.100	50.000	0.2
Endosulfan I	5.245	5.147	5.347	49.820	50.000	-0.4
Endosulfan II	6.080	5.981	6.181	49.230	50.000	-1.5
Endosulfan sulfate	6.482	6.383	6.583	48.600	50.000	-2.8
Endrin	5.789	5.689	5.889	50.440	50.000	0.9
Endrin aldehyde	6.258	6.159	6.359	47.730	50.000	-4.5
Endrin ketone	6.991	6.892	7.092	49.210	50.000	-1.6
gamma-BHC (Lindane)	3.728	3.630	3.830	51.080	50.000	2.2
gamma-Chlordane	5.124	5.026	5.226	49.420	50.000	-1.2
Heptachlor	4.083	3.983	4.183	49.000	50.000	-2.0
Heptachlor epoxide	4.871	4.773	4.973	50.730	50.000	1.5
Methoxychlor	6.754	6.655	6.855	51.030	50.000	2.1
Tetrachloro-m-xylene	2.881	2.780	2.980	53.140	50.000	6.3

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089277.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 16:50
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:48:02 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	148.0E6	901.0E6	51.883	53.144
28)	SA Decachlor...	9.073	8.072	191.2E6	943.0E6	48.693	47.701
Target Compounds							
2)	A alpha-BHC	4.000	3.394	304.7E6	1380.2E6	52.275	51.505
3)	MA gamma-BHC...	4.331	3.728	289.1E6	1264.7E6	51.802	51.083m
4)	MA Heptachlor	4.930	4.083	273.7E6	1226.8E6	50.196	48.995
5)	MB Aldrin	5.272	4.369	274.6E6	1238.1E6	51.685	51.126
6)	B beta-BHC	4.516	4.026	110.1E6	531.8E6	49.903	49.491
7)	B delta-BHC	4.765	4.263	277.1E6	1259.5E6	53.989	50.637
8)	B Heptachlo...	5.691	4.871	241.0E6	1111.0E6	49.950	50.733m
9)	A Endosulfan I	6.075	5.245	228.0E6	1038.3E6	50.576	49.818m
10)	B gamma-Chl...	5.947	5.124	244.0E6	1182.2E6	50.961	49.416m
11)	B alpha-Chl...	6.028	5.189	243.1E6	1139.0E6	50.237	49.746m
12)	B 4,4'-DDE	6.196	5.375	214.7E6	1151.6E6	49.239	50.252
13)	MA Dieldrin	6.348	5.511	240.8E6	1162.8E6	50.194	50.101m
14)	MA Endrin	6.574	5.789	194.8E6	1095.3E6	47.216	50.441
15)	B Endosulfa...	6.787	6.080	199.6E6	994.4E6	49.447	49.234
16)	A 4,4'-DDD	6.706	5.930	177.8E6	919.2E6	51.810	48.050
17)	MA 4,4'-DDT	7.021	6.183	167.3E6	894.8E6	44.174	44.124
18)	B Endrin al...	6.916	6.258	148.1E6	730.8E6	48.353	47.728
19)	B Endosulfa...	7.150	6.482	185.4E6	953.6E6	48.310	48.600
20)	A Methoxychlor	7.493	6.754	101.4E6	549.5E6	50.180	51.028
21)	B Endrin ke...	7.630	6.991	200.1E6	1063.0E6	49.127	49.211
22)	Mirex	8.114	7.184	144.8E6	791.1E6	47.644	47.430

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089277.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 16:50
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

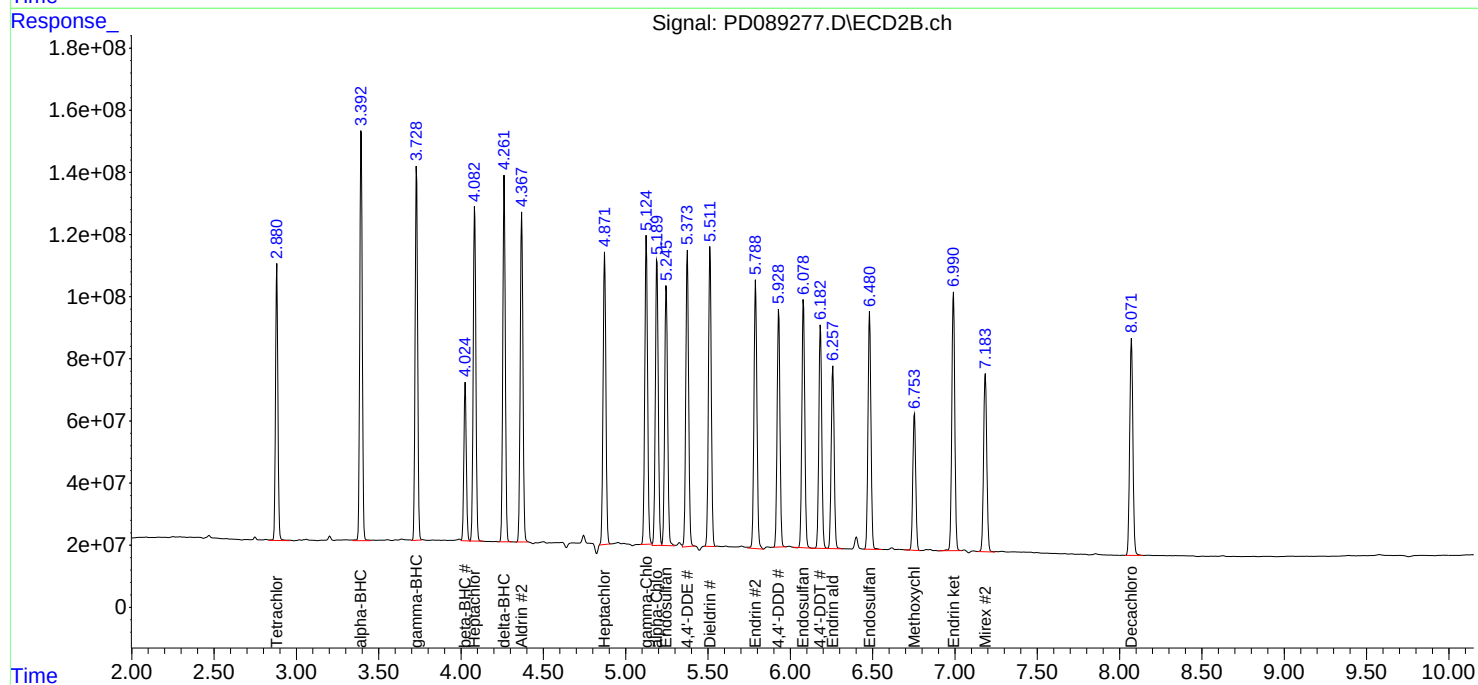
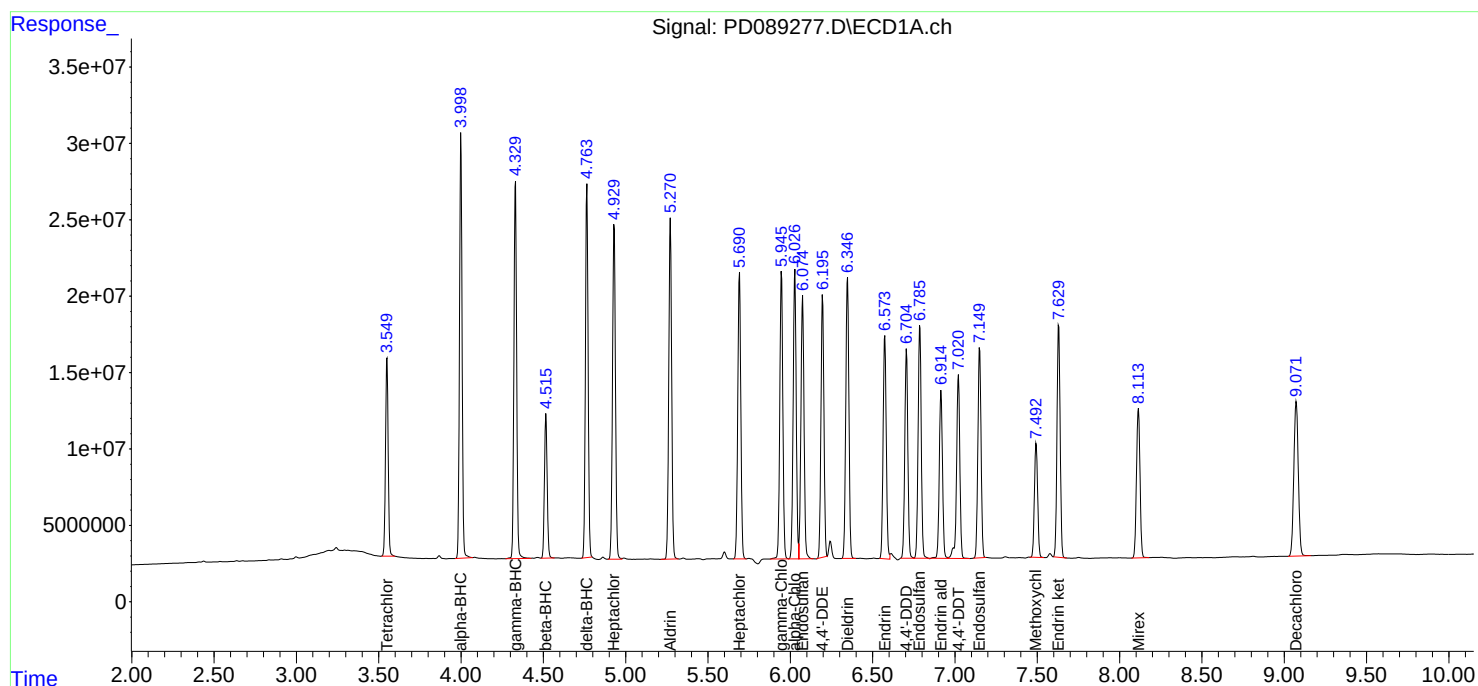
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:48:02 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 19:34

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.07	8.97	9.17	0.00
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	4.00	3.90	4.10	0.00
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.76	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.93	4.83	5.03	0.00
Aldrin	5.27	5.27	5.17	5.37	0.00
Heptachlor epoxide	5.69	5.69	5.59	5.79	0.00
Endosulfan I	6.08	6.08	5.98	6.18	0.01
Dieldrin	6.35	6.35	6.25	6.45	0.00
4,4'-DDE	6.20	6.20	6.10	6.30	0.00
Endrin	6.58	6.58	6.48	6.68	0.01
Endosulfan II	6.79	6.79	6.69	6.89	0.00
4,4'-DDD	6.71	6.71	6.61	6.81	0.01
Endosulfan sulfate	7.15	7.15	7.05	7.25	0.00
4,4'-DDT	7.02	7.02	6.92	7.12	0.00
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.92	6.82	7.02	0.01
alpha-Chlordane	6.03	6.03	5.93	6.13	0.00
gamma-Chlordane	5.95	5.95	5.85	6.05	0.00



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 19:34

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.00
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.38	5.38	5.28	5.48	0.00
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.18	6.18	6.08	6.28	0.00
Methoxychlor	6.75	6.76	6.66	6.86	0.01
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.12	5.13	5.03	5.23	0.01



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL03 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089289.D **Time Analyzed:** 19:34

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.705	6.605	6.805	50.410	50.000	0.8
4,4'-DDE	6.196	6.096	6.296	48.040	50.000	-3.9
4,4'-DDT	7.021	6.921	7.121	42.110	50.000	-15.8
Aldrin	5.271	5.171	5.371	50.960	50.000	1.9
alpha-BHC	3.999	3.899	4.099	51.940	50.000	3.9
alpha-Chlordane	6.027	5.927	6.127	49.270	50.000	-1.5
beta-BHC	4.516	4.415	4.615	49.360	50.000	-1.3
Decachlorobiphenyl	9.072	8.972	9.172	46.970	50.000	-6.1
delta-BHC	4.764	4.664	4.864	53.460	50.000	6.9
Dieldrin	6.347	6.247	6.447	49.220	50.000	-1.6
Endosulfan I	6.075	5.975	6.175	49.550	50.000	-0.9
Endosulfan II	6.786	6.686	6.886	47.770	50.000	-4.5
Endosulfan sulfate	7.149	7.049	7.249	46.540	50.000	-6.9
Endrin	6.575	6.475	6.675	46.290	50.000	-7.4
Endrin aldehyde	6.915	6.815	7.015	47.210	50.000	-5.6
Endrin ketone	7.630	7.530	7.730	47.690	50.000	-4.6
gamma-BHC (Lindane)	4.331	4.230	4.430	51.280	50.000	2.6
gamma-Chlordane	5.946	5.846	6.046	49.990	50.000	0.0
Heptachlor	4.930	4.829	5.029	49.310	50.000	-1.4
Heptachlor epoxide	5.692	5.591	5.791	48.920	50.000	-2.2
Methoxychlor	7.493	7.393	7.593	48.080	50.000	-3.8
Tetrachloro-m-xylene	3.550	3.450	3.650	51.730	50.000	3.5



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL03 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089289.D **Time Analyzed:** 19:34

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.930	5.830	6.030	46.490	50.000	-7.0
4,4'-DDE	5.375	5.275	5.475	48.790	50.000	-2.4
4,4'-DDT	6.183	6.084	6.284	41.790	50.000	-16.4
Aldrin	4.369	4.269	4.469	49.870	50.000	-0.3
alpha-BHC	3.393	3.293	3.493	50.900	50.000	1.8
alpha-Chlordane	5.190	5.091	5.291	50.080	50.000	0.2
beta-BHC	4.026	3.926	4.126	48.780	50.000	-2.4
Decachlorobiphenyl	8.072	7.972	8.172	44.010	50.000	-12.0
delta-BHC	4.262	4.162	4.362	49.450	50.000	-1.1
Dieldrin	5.511	5.413	5.613	48.580	50.000	-2.8
Endosulfan I	5.247	5.147	5.347	50.020	50.000	0.0
Endosulfan II	6.080	5.981	6.181	47.250	50.000	-5.5
Endosulfan sulfate	6.482	6.383	6.583	46.320	50.000	-7.4
Endrin	5.789	5.689	5.889	48.800	50.000	-2.4
Endrin aldehyde	6.259	6.159	6.359	45.640	50.000	-8.7
Endrin ketone	6.991	6.892	7.092	46.530	50.000	-6.9
gamma-BHC (Lindane)	3.730	3.630	3.830	49.230	50.000	-1.5
gamma-Chlordane	5.124	5.026	5.226	48.400	50.000	-3.2
Heptachlor	4.083	3.983	4.183	47.790	50.000	-4.4
Heptachlor epoxide	4.871	4.773	4.973	49.490	50.000	-1.0
Methoxychlor	6.753	6.655	6.855	48.100	50.000	-3.8
Tetrachloro-m-xylene	2.881	2.780	2.980	52.340	50.000	4.7

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089289.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 19:34
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:50:47 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	147.6E6	887.4E6	51.727	52.342
2)	SA Decachlor...	9.072	8.072	184.5E6	870.1E6	46.969	44.012
Target Compounds							
2)	A alpha-BHC	3.999	3.393	302.8E6	1363.9E6	51.945	50.897
3)	MA gamma-BHC...	4.331	3.730	286.3E6	1218.8E6	51.284	49.229
4)	MA Heptachlor	4.930	4.083	268.8E6	1196.6E6	49.306	47.792
5)	MB Aldrin	5.271	4.369	270.8E6	1207.8E6	50.962	49.874
6)	B beta-BHC	4.516	4.026	108.9E6	524.2E6	49.361	48.777
7)	B delta-BHC	4.764	4.262	274.4E6	1229.9E6	53.464	49.447
8)	B Heptachlo...	5.692	4.871	236.0E6	1083.7E6	48.922	49.488m
9)	A Endosulfan I	6.075	5.247	223.4E6	1042.5E6	49.554	50.018
10)	B gamma-Chl...	5.946	5.124	239.4E6	1158.0E6	49.988	48.405m
11)	B alpha-Chl...	6.027	5.190	238.4E6	1146.7E6	49.274	50.080
12)	B 4,4'-DDE	6.196	5.375	209.5E6	1118.1E6	48.037	48.792
13)	MA Dieldrin	6.347	5.511	236.1E6	1127.4E6	49.220	48.577m
14)	MA Endrin	6.575	5.789	191.0E6	1059.5E6	46.291	48.795
15)	B Endosulfa...	6.786	6.080	192.8E6	954.4E6	47.772	47.253
16)	A 4,4'-DDD	6.705	5.930	173.0E6	889.3E6	50.409	46.487
17)	MA 4,4'-DDT	7.021	6.183	159.5E6	847.5E6	42.105	41.791
18)	B Endrin al...	6.915	6.259	144.7E6	698.9E6	47.211	45.644
19)	B Endosulfa...	7.149	6.482	178.6E6	908.8E6	46.536	46.315
20)	A Methoxychlor	7.493	6.753	97144839	518.0E6	48.077	48.105
21)	B Endrin ke...	7.630	6.991	194.2E6	1005.1E6	47.688	46.528
22)	Mirex	8.113	7.185	137.8E6	767.5E6	45.326	46.019

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089289.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 19:34
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

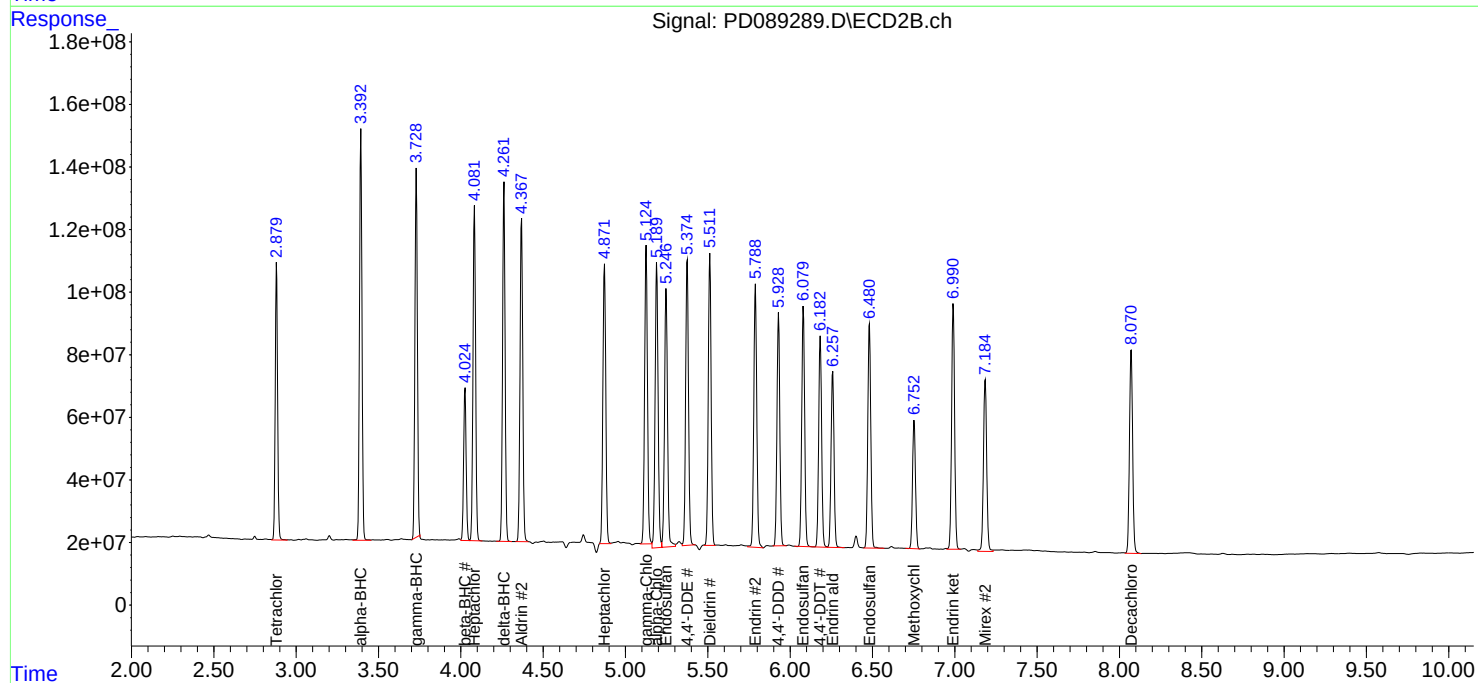
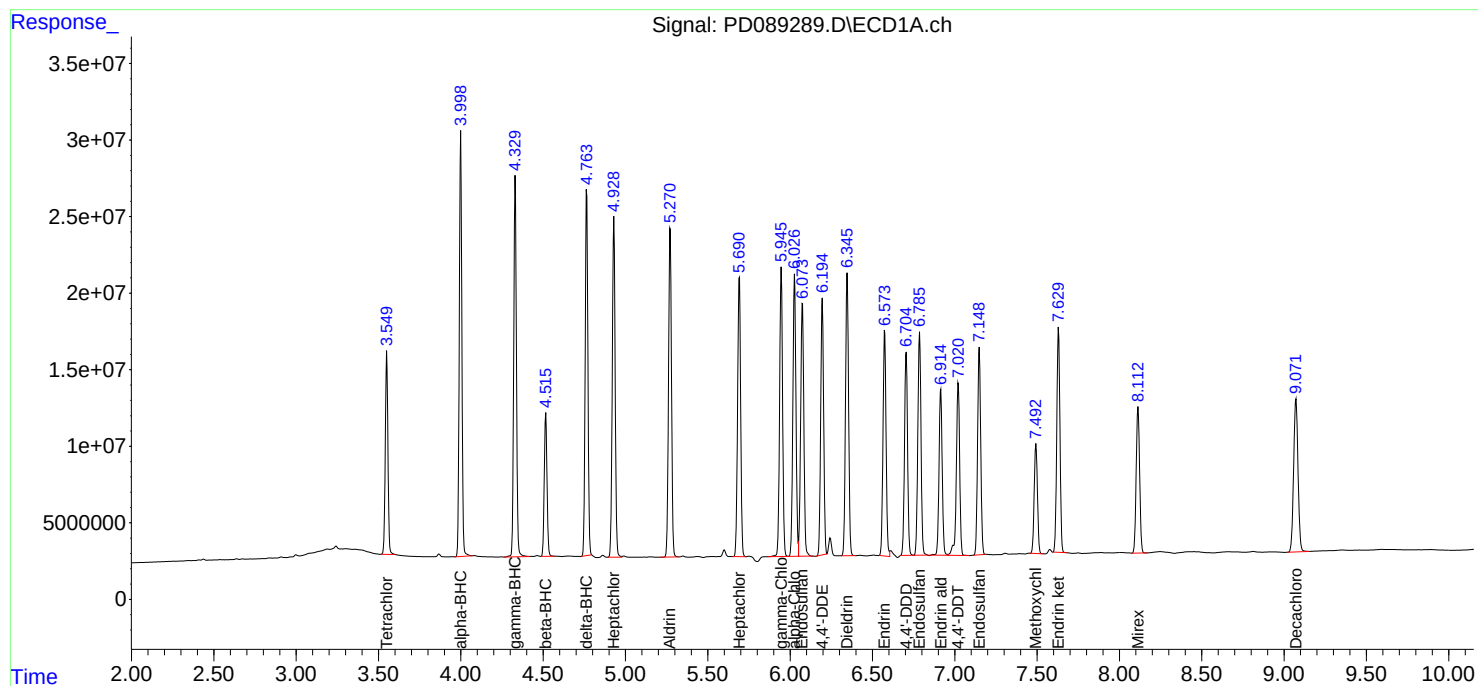
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:50:47 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 22:59

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.07	8.97	9.17	0.00
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	4.00	3.90	4.10	0.00
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.76	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.93	4.83	5.03	0.00
Aldrin	5.27	5.27	5.17	5.37	0.00
Heptachlor epoxide	5.69	5.69	5.59	5.79	0.00
Endosulfan I	6.07	6.08	5.98	6.18	0.01
Dieldrin	6.35	6.35	6.25	6.45	0.00
4,4'-DDE	6.20	6.20	6.10	6.30	0.00
Endrin	6.57	6.58	6.48	6.68	0.01
Endosulfan II	6.79	6.79	6.69	6.89	0.00
4,4'-DDD	6.71	6.71	6.61	6.81	0.01
Endosulfan sulfate	7.15	7.15	7.05	7.25	0.00
4,4'-DDT	7.02	7.02	6.92	7.12	0.00
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.92	6.82	7.02	0.01
alpha-Chlordane	6.03	6.03	5.93	6.13	0.00
gamma-Chlordane	5.95	5.95	5.85	6.05	0.00



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/01/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 22:59

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.00
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.38	5.38	5.28	5.48	0.00
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.18	6.18	6.08	6.28	0.00
Methoxychlor	6.75	6.76	6.66	6.86	0.01
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.13	5.13	5.03	5.23	0.00



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL04 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089299.D **Time Analyzed:** 22:59

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.705	6.605	6.805	49.240	50.000	-1.5
4,4'-DDE	6.196	6.096	6.296	47.280	50.000	-5.4
4,4'-DDT	7.021	6.921	7.121	40.210	50.000	-19.6
Aldrin	5.271	5.171	5.371	50.470	50.000	0.9
alpha-BHC	4.000	3.899	4.099	51.690	50.000	3.4
alpha-Chlordane	6.027	5.927	6.127	48.830	50.000	-2.3
beta-BHC	4.516	4.415	4.615	49.400	50.000	-1.2
Decachlorobiphenyl	9.073	8.972	9.172	45.290	50.000	-9.4
delta-BHC	4.764	4.664	4.864	53.160	50.000	6.3
Dieldrin	6.347	6.247	6.447	48.690	50.000	-2.6
Endosulfan I	6.074	5.975	6.175	48.840	50.000	-2.3
Endosulfan II	6.786	6.686	6.886	46.800	50.000	-6.4
Endosulfan sulfate	7.150	7.049	7.249	45.830	50.000	-8.3
Endrin	6.574	6.475	6.675	44.690	50.000	-10.6
Endrin aldehyde	6.915	6.815	7.015	45.270	50.000	-9.5
Endrin ketone	7.630	7.530	7.730	46.190	50.000	-7.6
gamma-BHC (Lindane)	4.330	4.230	4.430	51.150	50.000	2.3
gamma-Chlordane	5.946	5.846	6.046	49.870	50.000	-0.3
Heptachlor	4.930	4.829	5.029	48.000	50.000	-4.0
Heptachlor epoxide	5.690	5.591	5.791	48.490	50.000	-3.0
Methoxychlor	7.493	7.393	7.593	45.780	50.000	-8.4
Tetrachloro-m-xylene	3.550	3.450	3.650	51.450	50.000	2.9



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL04 **Date Analyzed:** 07/01/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089299.D **Time Analyzed:** 22:59

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.929	5.830	6.030	48.080	50.000	-3.8
4,4'-DDE	5.375	5.275	5.475	48.280	50.000	-3.4
4,4'-DDT	6.182	6.084	6.284	40.910	50.000	-18.2
Aldrin	4.368	4.269	4.469	50.130	50.000	0.3
alpha-BHC	3.393	3.293	3.493	50.710	50.000	1.4
alpha-Chlordane	5.189	5.091	5.291	48.800	50.000	-2.4
beta-BHC	4.025	3.926	4.126	48.630	50.000	-2.7
Decachlorobiphenyl	8.071	7.972	8.172	41.860	50.000	-16.3
delta-BHC	4.262	4.162	4.362	49.480	50.000	-1.0
Dieldrin	5.511	5.413	5.613	48.470	50.000	-3.1
Endosulfan I	5.246	5.147	5.347	48.840	50.000	-2.3
Endosulfan II	6.080	5.981	6.181	47.030	50.000	-5.9
Endosulfan sulfate	6.481	6.383	6.583	45.240	50.000	-9.5
Endrin	5.789	5.689	5.889	48.700	50.000	-2.6
Endrin aldehyde	6.258	6.159	6.359	45.360	50.000	-9.3
Endrin ketone	6.990	6.892	7.092	45.230	50.000	-9.5
gamma-BHC (Lindane)	3.729	3.630	3.830	48.650	50.000	-2.7
gamma-Chlordane	5.125	5.026	5.226	49.370	50.000	-1.3
Heptachlor	4.082	3.983	4.183	47.260	50.000	-5.5
Heptachlor epoxide	4.872	4.773	4.973	51.730	50.000	3.5
Methoxychlor	6.753	6.655	6.855	46.030	50.000	-7.9
Tetrachloro-m-xylene	2.881	2.780	2.980	52.440	50.000	4.9

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089299.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 22:59
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:53:46 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	146.8E6	889.0E6	51.446	52.441
28)	SA Decachlor...	9.073	8.071	177.9E6	827.6E6	45.290	41.860
Target Compounds							
2)	A alpha-BHC	4.000	3.393	301.3E6	1358.8E6	51.687	50.705
3)	MA gamma-BHC...	4.330	3.729	285.5E6	1204.4E6	51.154	48.651
4)	MA Heptachlor	4.930	4.082	261.7E6	1183.4E6	48.000	47.262
5)	MB Aldrin	5.271	4.368	268.2E6	1214.1E6	50.472	50.132
6)	B beta-BHC	4.516	4.025	109.0E6	522.5E6	49.401	48.626
7)	B delta-BHC	4.764	4.262	272.9E6	1230.8E6	53.164	49.482
8)	B Heptachlo...	5.690	4.872	233.9E6	1132.8E6	48.492	51.733
9)	A Endosulfan I	6.074	5.246	220.1E6	1018.0E6	48.836	48.843
10)	B gamma-Chl...	5.946	5.125	238.8E6	1181.1E6	49.873	49.371
11)	B alpha-Chl...	6.027	5.189	236.3E6	1117.3E6	48.827	48.797
12)	B 4,4'-DDE	6.196	5.375	206.2E6	1106.5E6	47.276	48.284
13)	MA Dieldrin	6.347	5.511	233.6E6	1124.9E6	48.687	48.470m
14)	MA Endrin	6.574	5.789	184.4E6	1057.5E6	44.687	48.703
15)	B Endosulfa...	6.786	6.080	188.9E6	949.8E6	46.801	47.028
16)	A 4,4'-DDD	6.705	5.929	169.0E6	919.8E6	49.237	48.081
17)	MA 4,4'-DDT	7.021	6.182	152.3E6	829.6E6	40.212	40.907
18)	B Endrin al...	6.915	6.258	138.7E6	694.5E6	45.270	45.358
19)	B Endosulfa...	7.150	6.481	175.9E6	887.6E6	45.832	45.237
20)	A Methoxychlor	7.493	6.753	92501875	495.6E6	45.779	46.027
21)	B Endrin ke...	7.630	6.990	188.1E6	977.2E6	46.186	45.235
22)	Mirex	8.114	7.184	137.0E6	721.0E6	45.048	43.233

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089299.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 22:59
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

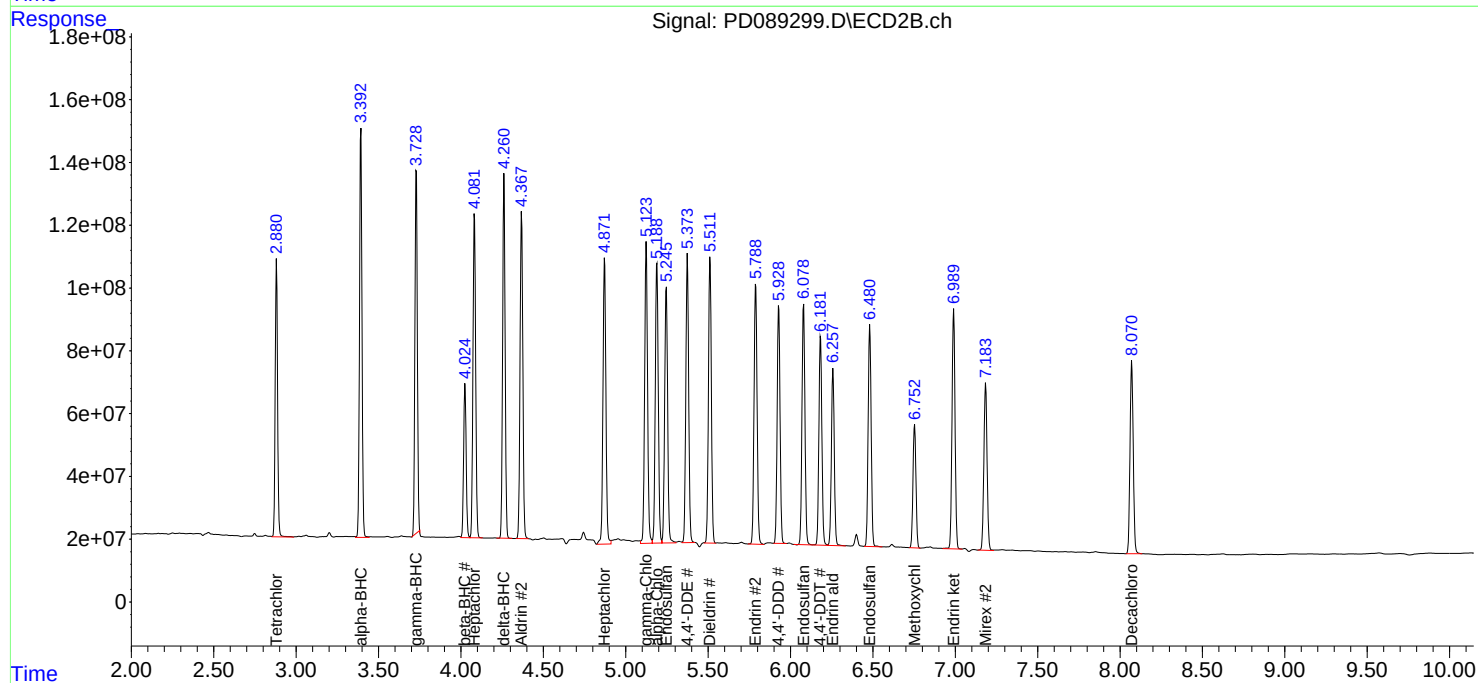
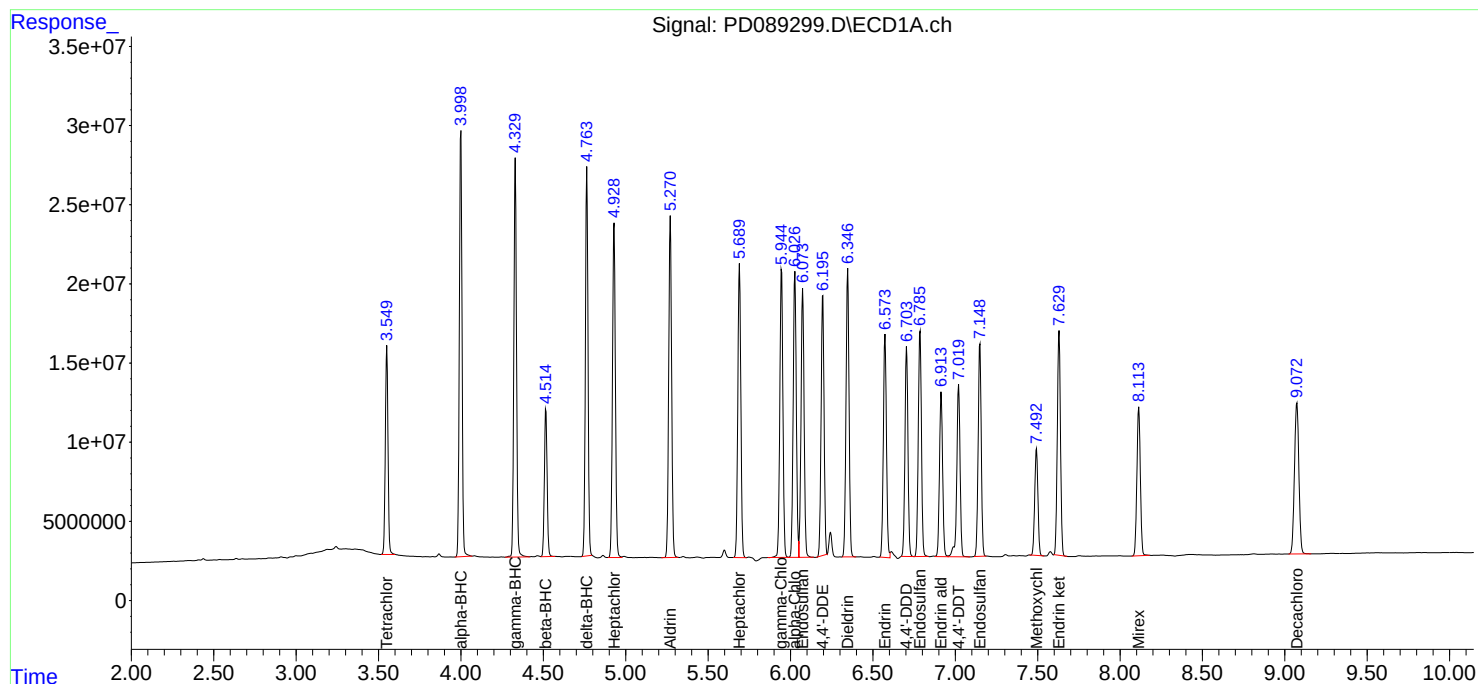
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:53:46 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/03/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 09:46

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.07	8.97	9.17	0.00
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	4.00	3.90	4.10	0.00
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.77	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.93	4.83	5.03	0.00
Aldrin	5.27	5.27	5.17	5.37	0.00
Heptachlor epoxide	5.69	5.69	5.59	5.79	0.00
Endosulfan I	6.08	6.08	5.98	6.18	0.01
Dieldrin	6.35	6.35	6.25	6.45	0.00
4,4'-DDE	6.20	6.20	6.10	6.30	0.00
Endrin	6.57	6.58	6.48	6.68	0.01
Endosulfan II	6.79	6.79	6.69	6.89	0.00
4,4'-DDD	6.71	6.71	6.61	6.81	0.01
Endosulfan sulfate	7.15	7.15	7.05	7.25	0.00
4,4'-DDT	7.02	7.02	6.92	7.12	0.00
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.92	6.82	7.02	0.01
alpha-Chlordane	6.03	6.03	5.93	6.13	0.00
gamma-Chlordane	5.95	5.95	5.85	6.05	0.00



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Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/03/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 09:46

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.00
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.37	5.38	5.28	5.48	0.01
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.18	6.18	6.08	6.28	0.00
Methoxychlor	6.75	6.76	6.66	6.86	0.01
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.12	5.13	5.03	5.23	0.01



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL05 **Date Analyzed:** 07/03/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089328.D **Time Analyzed:** 09:46

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.705	6.605	6.805	54.350	50.000	8.7
4,4'-DDE	6.196	6.096	6.296	51.990	50.000	4.0
4,4'-DDT	7.021	6.921	7.121	43.940	50.000	-12.1
Aldrin	5.271	5.171	5.371	55.520	50.000	11.0
alpha-BHC	4.000	3.899	4.099	56.340	50.000	12.7
alpha-Chlordane	6.027	5.927	6.127	54.210	50.000	8.4
beta-BHC	4.516	4.415	4.615	53.430	50.000	6.9
Decachlorobiphenyl	9.073	8.972	9.172	46.920	50.000	-6.2
delta-BHC	4.765	4.664	4.864	58.070	50.000	16.1
Dieldrin	6.347	6.247	6.447	54.110	50.000	8.2
Endosulfan I	6.075	5.975	6.175	54.180	50.000	8.4
Endosulfan II	6.786	6.686	6.886	51.490	50.000	3.0
Endosulfan sulfate	7.150	7.049	7.249	49.420	50.000	-1.2
Endrin	6.574	6.475	6.675	49.350	50.000	-1.3
Endrin aldehyde	6.915	6.815	7.015	49.610	50.000	-0.8
Endrin ketone	7.630	7.530	7.730	49.570	50.000	-0.9
gamma-BHC (Lindane)	4.331	4.230	4.430	55.920	50.000	11.8
gamma-Chlordane	5.946	5.846	6.046	55.190	50.000	10.4
Heptachlor	4.930	4.829	5.029	52.740	50.000	5.5
Heptachlor epoxide	5.691	5.591	5.791	53.220	50.000	6.4
Methoxychlor	7.493	7.393	7.593	49.780	50.000	-0.4
Tetrachloro-m-xylene	3.550	3.450	3.650	55.630	50.000	11.3



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CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL05 **Date Analyzed:** 07/03/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089328.D **Time Analyzed:** 09:46

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.929	5.830	6.030	52.750	50.000	5.5
4,4'-DDE	5.374	5.275	5.475	52.830	50.000	5.7
4,4'-DDT	6.183	6.084	6.284	44.690	50.000	-10.6
Aldrin	4.368	4.269	4.469	54.130	50.000	8.3
alpha-BHC	3.393	3.293	3.493	55.810	50.000	11.6
alpha-Chlordane	5.189	5.091	5.291	53.210	50.000	6.4
beta-BHC	4.025	3.926	4.126	51.100	50.000	2.2
Decachlorobiphenyl	8.072	7.972	8.172	44.450	50.000	-11.1
delta-BHC	4.261	4.162	4.362	53.310	50.000	6.6
Dieldrin	5.511	5.413	5.613	53.310	50.000	6.6
Endosulfan I	5.246	5.147	5.347	53.080	50.000	6.2
Endosulfan II	6.080	5.981	6.181	51.600	50.000	3.2
Endosulfan sulfate	6.482	6.383	6.583	49.050	50.000	-1.9
Endrin	5.789	5.689	5.889	54.650	50.000	9.3
Endrin aldehyde	6.258	6.159	6.359	49.950	50.000	-0.1
Endrin ketone	6.990	6.892	7.092	48.890	50.000	-2.2
gamma-BHC (Lindane)	3.728	3.630	3.830	54.280	50.000	8.6
gamma-Chlordane	5.123	5.026	5.226	52.410	50.000	4.8
Heptachlor	4.082	3.983	4.183	50.420	50.000	0.8
Heptachlor epoxide	4.872	4.773	4.973	56.110	50.000	12.2
Methoxychlor	6.754	6.655	6.855	49.550	50.000	-0.9
Tetrachloro-m-xylene	2.881	2.780	2.980	57.330	50.000	14.7

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089328.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 09:46
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:17:08 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	158.7E6	971.8E6	55.633	57.326
2)	SA Decachlor...	9.073	8.072	184.3E6	878.8E6	46.924	44.451
Target Compounds							
2)	A alpha-BHC	4.000	3.393	328.4E6	1495.5E6	56.340	55.809
3)	MA gamma-BHC...	4.331	3.728	312.1E6	1343.7E6	55.916	54.276m
4)	MA Heptachlor	4.930	4.082	287.5E6	1262.3E6	52.736	50.416
5)	MB Aldrin	5.271	4.368	295.0E6	1310.9E6	55.520	54.132
6)	B beta-BHC	4.516	4.025	117.9E6	549.1E6	53.431	51.099
7)	B delta-BHC	4.765	4.261	298.1E6	1325.9E6	58.071	53.307
8)	B Heptachlo...	5.691	4.872	256.7E6	1228.7E6	53.216	56.111
9)	A Endosulfan I	6.075	5.246	244.2E6	1106.3E6	54.183	53.078
10)	B gamma-Chl...	5.946	5.123	264.2E6	1253.8E6	55.187	52.409m
11)	B alpha-Chl...	6.027	5.189	262.3E6	1218.3E6	54.214	53.210
12)	B 4,4'-DDE	6.196	5.374	226.7E6	1210.6E6	51.990	52.826
13)	MA Dieldrin	6.347	5.511	259.6E6	1237.2E6	54.113	53.308m
14)	MA Endrin	6.574	5.789	203.6E6	1186.6E6	49.348	54.648
15)	B Endosulfa...	6.786	6.080	207.8E6	1042.1E6	51.485	51.597
16)	A 4,4'-DDD	6.705	5.929	186.5E6	1009.1E6	54.347	52.753
17)	MA 4,4'-DDT	7.021	6.183	166.4E6	906.3E6	43.936	44.691
18)	B Endrin al...	6.915	6.258	152.0E6	764.8E6	49.611	49.948
19)	B Endosulfa...	7.150	6.482	189.6E6	962.3E6	49.419	49.046
20)	A Methoxychlor	7.493	6.754	100.6E6	533.6E6	49.778	49.551
21)	B Endrin ke...	7.630	6.990	201.9E6	1056.2E6	49.575	48.893
22)	Mirex	8.114	7.185	144.4E6	787.7E6	47.482	47.230

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089328.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 09:46
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

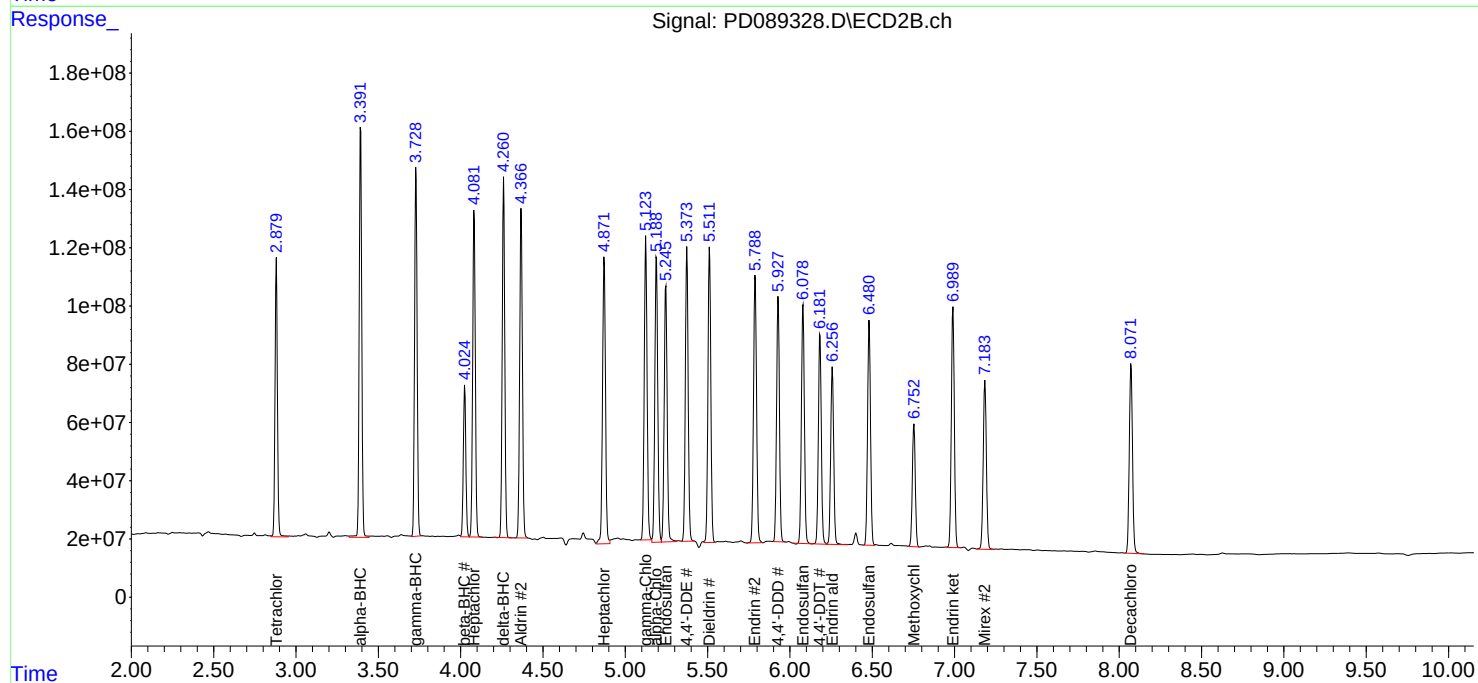
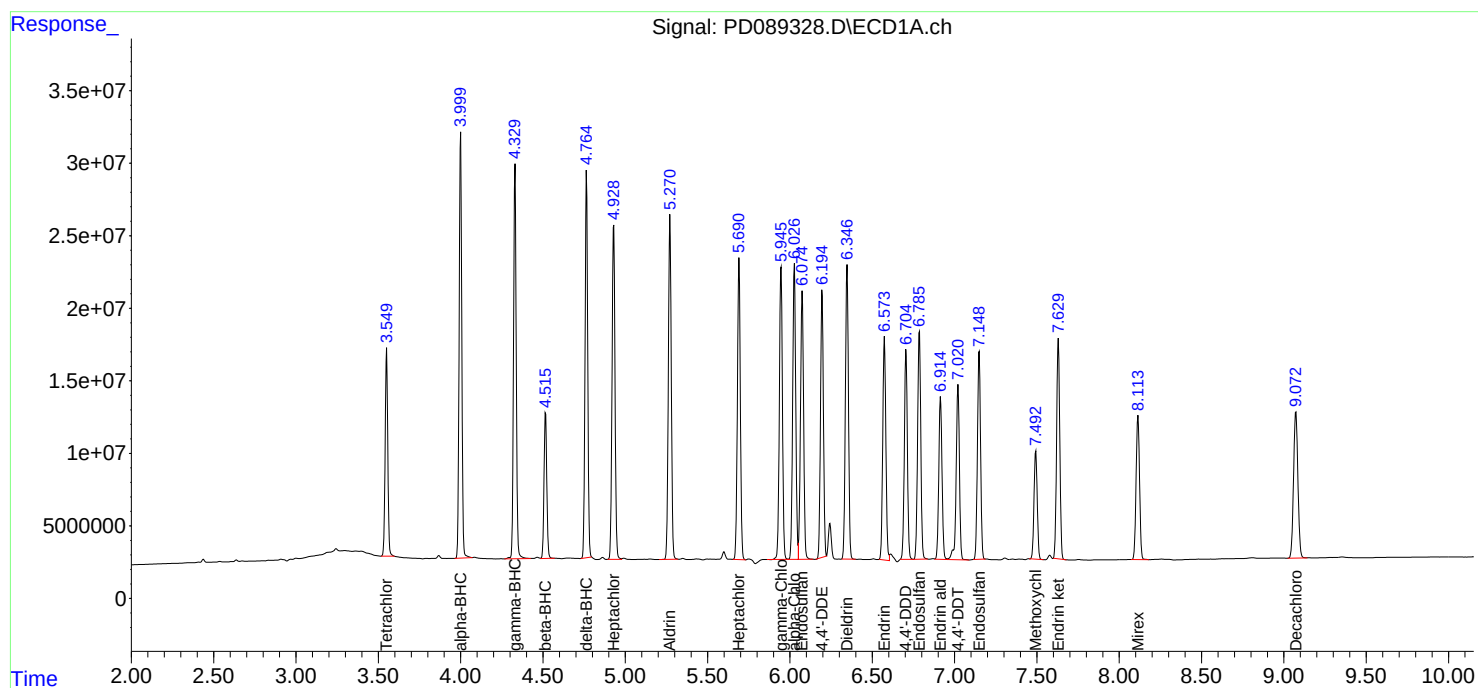
APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:17:08 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





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Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/03/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 16:02

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR1 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.07	8.97	9.17	0.00
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	4.00	3.90	4.10	0.00
beta-BHC	4.52	4.52	4.42	4.62	0.00
delta-BHC	4.77	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.93	4.83	5.03	0.00
Aldrin	5.27	5.27	5.17	5.37	0.00
Heptachlor epoxide	5.69	5.69	5.59	5.79	0.00
Endosulfan I	6.08	6.08	5.98	6.18	0.01
Dieldrin	6.35	6.35	6.25	6.45	0.00
4,4'-DDE	6.20	6.20	6.10	6.30	0.00
Endrin	6.58	6.58	6.48	6.68	0.01
Endosulfan II	6.79	6.79	6.69	6.89	0.00
4,4'-DDD	6.71	6.71	6.61	6.81	0.00
Endosulfan sulfate	7.15	7.15	7.05	7.25	0.00
4,4'-DDT	7.02	7.02	6.92	7.12	0.00
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.92	6.82	7.02	0.00
alpha-Chlordane	6.03	6.03	5.93	6.13	0.00
gamma-Chlordane	5.95	5.95	5.85	6.05	0.00



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Continuing Calib Date: 07/03/2025

Initial Calibration Date(s): 06/17/2025

06/17/2025

Continuing Calib Time: 16:02

Initial Calibration Time(s): 15:52

16:47

GC Column: ZB-MR2 ID: 0.32 (mm)

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.07	7.97	8.17	0.00
Tetrachloro-m-xylene	2.88	2.88	2.78	2.98	0.00
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.03	4.03	3.93	4.13	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.73	3.73	3.63	3.83	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.00
Aldrin	4.37	4.37	4.27	4.47	0.00
Heptachlor epoxide	4.87	4.87	4.77	4.97	0.00
Endosulfan I	5.25	5.25	5.15	5.35	0.00
Dieldrin	5.51	5.51	5.41	5.61	0.00
4,4'-DDE	5.38	5.38	5.28	5.48	0.00
Endrin	5.79	5.79	5.69	5.89	0.00
Endosulfan II	6.08	6.08	5.98	6.18	0.00
4,4'-DDD	5.93	5.93	5.83	6.03	0.00
Endosulfan sulfate	6.48	6.48	6.38	6.58	0.00
4,4'-DDT	6.18	6.18	6.08	6.28	0.00
Methoxychlor	6.75	6.76	6.66	6.86	0.01
Endrin ketone	6.99	6.99	6.89	7.09	0.00
Endrin aldehyde	6.26	6.26	6.16	6.36	0.00
alpha-Chlordane	5.19	5.19	5.09	5.29	0.00
gamma-Chlordane	5.13	5.13	5.03	5.23	0.00



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group Contract: CAMP02

Lab Code: CHEM SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No.: CCAL06 Date Analyzed: 07/03/2025

Lab Sample No.: PSTDCCC050 Data File : PD089338.D Time Analyzed: 16:02

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.706	6.605	6.805	54.570	50.000	9.1
4,4'-DDE	6.196	6.096	6.296	50.560	50.000	1.1
4,4'-DDT	7.022	6.921	7.121	44.490	50.000	-11.0
Aldrin	5.271	5.171	5.371	54.740	50.000	9.5
alpha-BHC	4.000	3.899	4.099	55.900	50.000	11.8
alpha-Chlordane	6.027	5.927	6.127	52.800	50.000	5.6
beta-BHC	4.516	4.415	4.615	52.760	50.000	5.5
Decachlorobiphenyl	9.074	8.972	9.172	46.330	50.000	-7.3
delta-BHC	4.765	4.664	4.864	57.140	50.000	14.3
Dieldrin	6.348	6.247	6.447	52.530	50.000	5.1
Endosulfan I	6.075	5.975	6.175	52.890	50.000	5.8
Endosulfan II	6.787	6.686	6.886	51.760	50.000	3.5
Endosulfan sulfate	7.151	7.049	7.249	50.190	50.000	0.4
Endrin	6.575	6.475	6.675	49.270	50.000	-1.5
Endrin aldehyde	6.916	6.815	7.015	50.280	50.000	0.6
Endrin ketone	7.632	7.530	7.730	50.190	50.000	0.4
gamma-BHC (Lindane)	4.331	4.230	4.430	55.390	50.000	10.8
gamma-Chlordane	5.947	5.846	6.046	53.600	50.000	7.2
Heptachlor	4.930	4.829	5.029	52.380	50.000	4.8
Heptachlor epoxide	5.691	5.591	5.791	52.180	50.000	4.4
Methoxychlor	7.494	7.393	7.593	50.700	50.000	1.4
Tetrachloro-m-xylene	3.551	3.450	3.650	55.250	50.000	10.5



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance Group **Contract:** CAMP02
Lab Code: CHEM **SDG NO.:** Q2458
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 06/17/2025 06/17/2025

Client Sample No.: CCAL06 **Date Analyzed:** 07/03/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD089338.D **Time Analyzed:** 16:02

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.929	5.830	6.030	52.350	50.000	4.7
4,4'-DDE	5.375	5.275	5.475	52.440	50.000	4.9
4,4'-DDT	6.184	6.084	6.284	45.480	50.000	-9.0
Aldrin	4.369	4.269	4.469	54.440	50.000	8.9
alpha-BHC	3.393	3.293	3.493	56.730	50.000	13.5
alpha-Chlordane	5.190	5.091	5.291	52.940	50.000	5.9
beta-BHC	4.025	3.926	4.126	51.210	50.000	2.4
Decachlorobiphenyl	8.072	7.972	8.172	46.160	50.000	-7.7
delta-BHC	4.262	4.162	4.362	54.020	50.000	8.0
Dieldrin	5.511	5.413	5.613	52.900	50.000	5.8
Endosulfan I	5.246	5.147	5.347	52.850	50.000	5.7
Endosulfan II	6.080	5.981	6.181	51.150	50.000	2.3
Endosulfan sulfate	6.482	6.383	6.583	50.040	50.000	0.1
Endrin	5.789	5.689	5.889	54.140	50.000	8.3
Endrin aldehyde	6.259	6.159	6.359	50.360	50.000	0.7
Endrin ketone	6.991	6.892	7.092	49.800	50.000	-0.4
gamma-BHC (Lindane)	3.728	3.630	3.830	55.020	50.000	10.0
gamma-Chlordane	5.125	5.026	5.226	53.820	50.000	7.6
Heptachlor	4.082	3.983	4.183	52.110	50.000	4.2
Heptachlor epoxide	4.872	4.773	4.973	56.280	50.000	12.6
Methoxychlor	6.754	6.655	6.855	52.330	50.000	4.7
Tetrachloro-m-xylene	2.881	2.780	2.980	57.920	50.000	15.8

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089338.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 16:02
 Operator : AR\AJ
 Sample : PSTDCCC050
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PSTDCCC050

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/07/2025
 Supervised By : mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:19:43 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.881	157.6E6	982.0E6	55.246	57.925
28)	SA Decachlor...	9.074	8.072	182.0E6	912.7E6	46.330	46.165
Target Compounds							
2)	A alpha-BHC	4.000	3.393	325.9E6	1520.2E6	55.904	56.730
3)	MA gamma-BHC...	4.331	3.728	309.2E6	1362.1E6	55.390	55.018m
4)	MA Heptachlor	4.930	4.082	285.6E6	1304.8E6	52.381	52.110
5)	MB Aldrin	5.271	4.369	290.8E6	1318.3E6	54.743	54.435
6)	B beta-BHC	4.516	4.025	116.4E6	550.4E6	52.760	51.214
7)	B delta-BHC	4.765	4.262	293.3E6	1343.7E6	57.139	54.022
8)	B Heptachlo...	5.691	4.872	251.7E6	1232.3E6	52.180	56.276
9)	A Endosulfan I	6.075	5.246	238.4E6	1101.6E6	52.893	52.854
10)	B gamma-Chl...	5.947	5.125	256.6E6	1287.6E6	53.600	53.823
11)	B alpha-Chl...	6.027	5.190	255.5E6	1212.2E6	52.805	52.941
12)	B 4,4'-DDE	6.196	5.375	220.5E6	1201.7E6	50.563	52.439
13)	MA Dieldrin	6.348	5.511	252.0E6	1227.7E6	52.534	52.899m
14)	MA Endrin	6.575	5.789	203.3E6	1175.5E6	49.269	54.136
15)	B Endosulfa...	6.787	6.080	208.9E6	1033.0E6	51.759	51.148
16)	A 4,4'-DDD	6.706	5.929	187.3E6	1001.5E6	54.571	52.352
17)	MA 4,4'-DDT	7.022	6.184	168.5E6	922.2E6	44.486	45.476
18)	B Endrin al...	6.916	6.259	154.1E6	771.1E6	50.279	50.364
19)	B Endosulfa...	7.151	6.482	192.6E6	981.9E6	50.194	50.043
20)	A Methoxychlor	7.494	6.754	102.4E6	563.5E6	50.696	52.335
21)	B Endrin ke...	7.632	6.991	204.4E6	1075.9E6	50.188	49.805
22)	Mirex	8.115	7.185	144.8E6	794.7E6	47.639	47.646

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089338.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 16:02
Operator : AR\AJ
Sample : PSTDCCC050
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PSTDCCC050

Manual Integrations

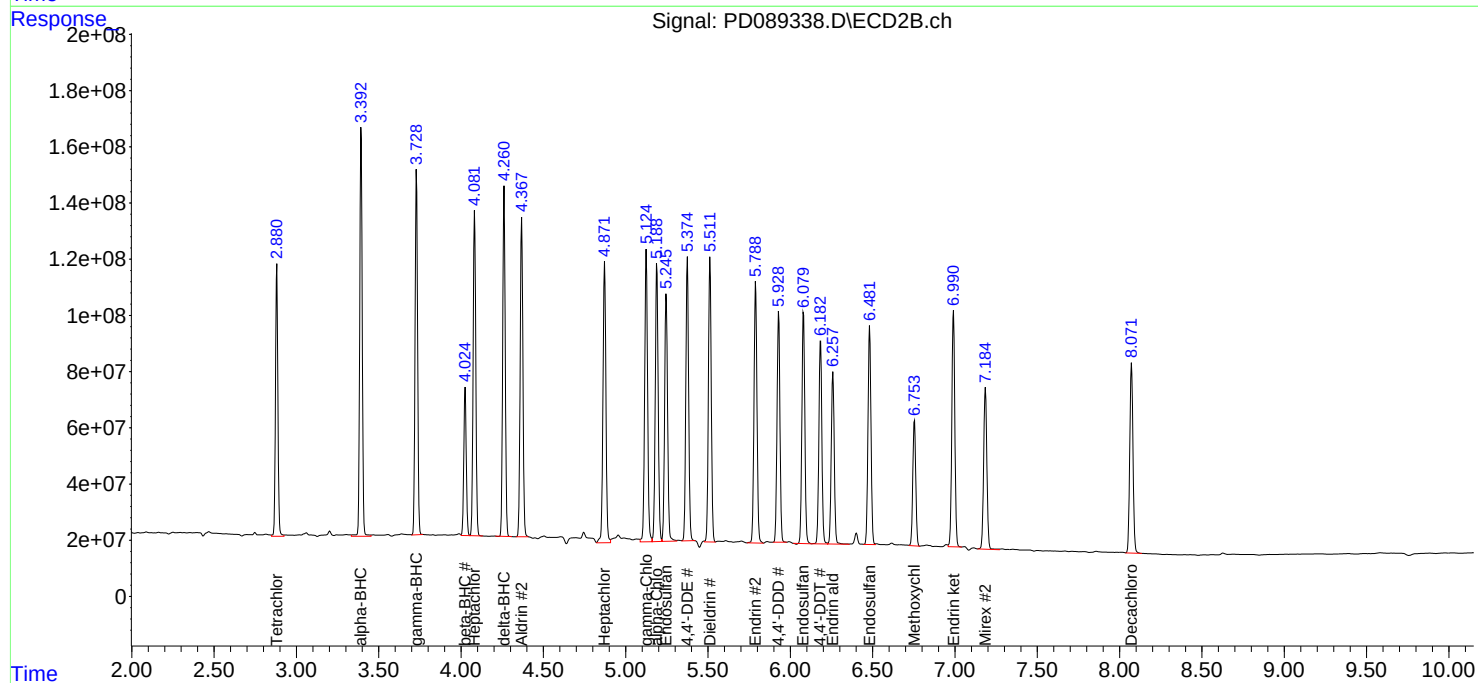
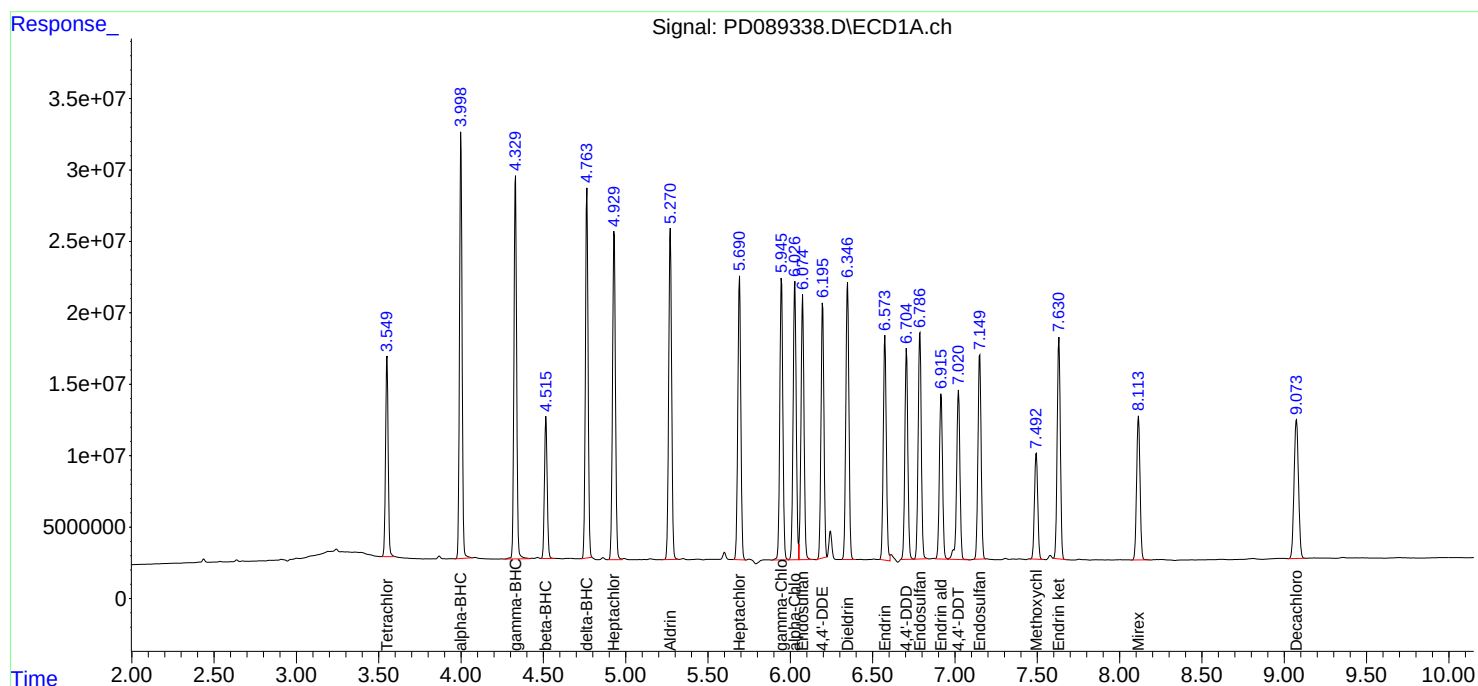
APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:19:43 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD088991.D Date Analyzed: 06/17/2025

Lab Sample No.(PEM): PEM Time Analyzed: 15:25

PEM COMPOUND	RT	RT WINDOW		CALC	NOM	%D
		FROM	TO	AMOUNT(ng)	AMOUNT(ng)	
Decachlorobiphenyl	9.071	8.970	9.170	21.440	20.000	7.2
Tetrachloro-m-xylene	3.550	3.500	3.600	19.860	20.000	-0.7
alpha-BHC	3.999	3.950	4.050	8.970	10.000	-10.3
beta-BHC	4.514	4.460	4.560	9.720	10.000	-2.8
gamma-BHC (Lindane)	4.330	4.280	4.380	9.320	10.000	-6.8
Endrin	6.574	6.500	6.640	50.260	50.000	0.5
4,4'-DDT	7.020	6.950	7.090	100.820	100.000	0.8
Methoxychlor	7.493	7.420	7.560	229.880	250.000	-8.0

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD088991.D Date Analyzed: 06/17/2025

Lab Sample No.(PEM): PEM Time Analyzed: 15:25

PEM COMPOUND	RT	RT WINDOW		CALC	NOM	%D
		FROM	TO	AMOUNT(ng)	AMOUNT(ng)	
Decachlorobiphenyl	8.072	7.970	8.170	21.290	20.000	6.5
Tetrachloro-m-xylene	2.880	2.830	2.930	20.440	20.000	2.2
alpha-BHC	3.393	3.340	3.440	10.570	10.000	5.7
beta-BHC	4.025	3.970	4.080	10.750	10.000	7.5
gamma-BHC (Lindane)	3.729	3.680	3.780	10.580	10.000	5.8
Endrin	5.789	5.720	5.860	50.490	50.000	1.0
4,4'-DDT	6.183	6.110	6.250	96.080	100.000	-3.9
Methoxychlor	6.754	6.680	6.820	198.860	250.000	-20.5

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
 Data File : PD088991.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 17 Jun 2025 15:25
 Operator : AR\AJ
 Sample : PEM
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PEM

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 06/18/2025
 Supervised By :mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jun 18 05:45:29 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.880	56676201	346.6E6	19.865	20.443
28)	SA Decachlor...	9.071	8.072	84217746	421.0E6	21.442	21.293
Target Compounds							
2)	A alpha-BHC	3.999	3.393	52304160	283.3E6	8.973	10.573
3)	MA gamma-BHC...	4.330	3.729	52036875	261.9E6	9.323	10.578
6)	B beta-BHC	4.514	4.025	21456971	115.6E6	9.722m	10.754
14)	MA Endrin	6.574	5.789	207.3E6	1096.4E6	50.256	50.492
16)	A 4,4'-DDD	6.701	5.931	603821	5602244	0.176m	0.293m#
17)	MA 4,4'-DDT	7.020	6.183	381.9E6	1948.4E6	100.821	96.080
18)	B Endrin al...	6.918	6.256	451112	9421924	0.147m	0.615m#
20)	A Methoxychlor	7.493	6.754	464.5E6	2141.4E6	229.876	198.863
21)	B Endrin ke...	7.627	6.989	1470511	10658868	0.361m	0.493m#

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088991.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 15:25
Operator : AR\AJ
Sample : PEM
Misc :
ALS Vial : 3 Sample Multiplier: 1

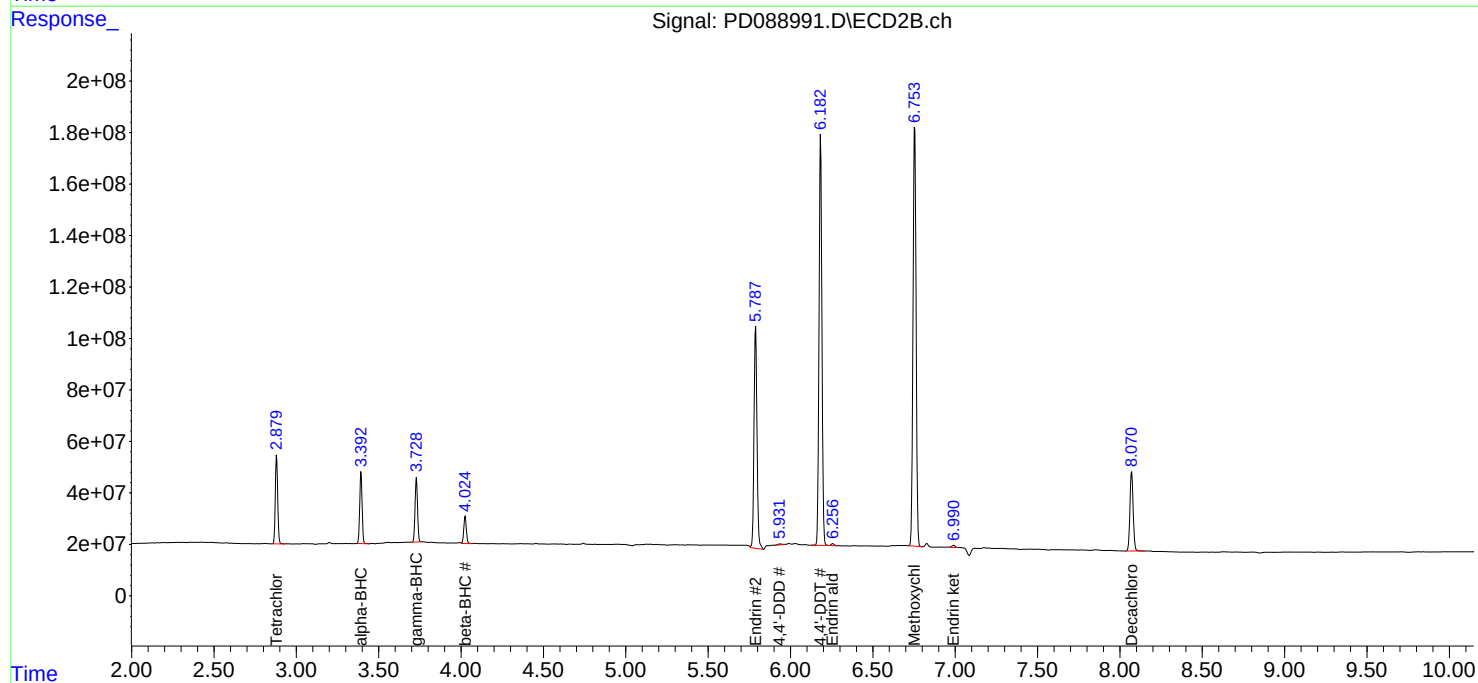
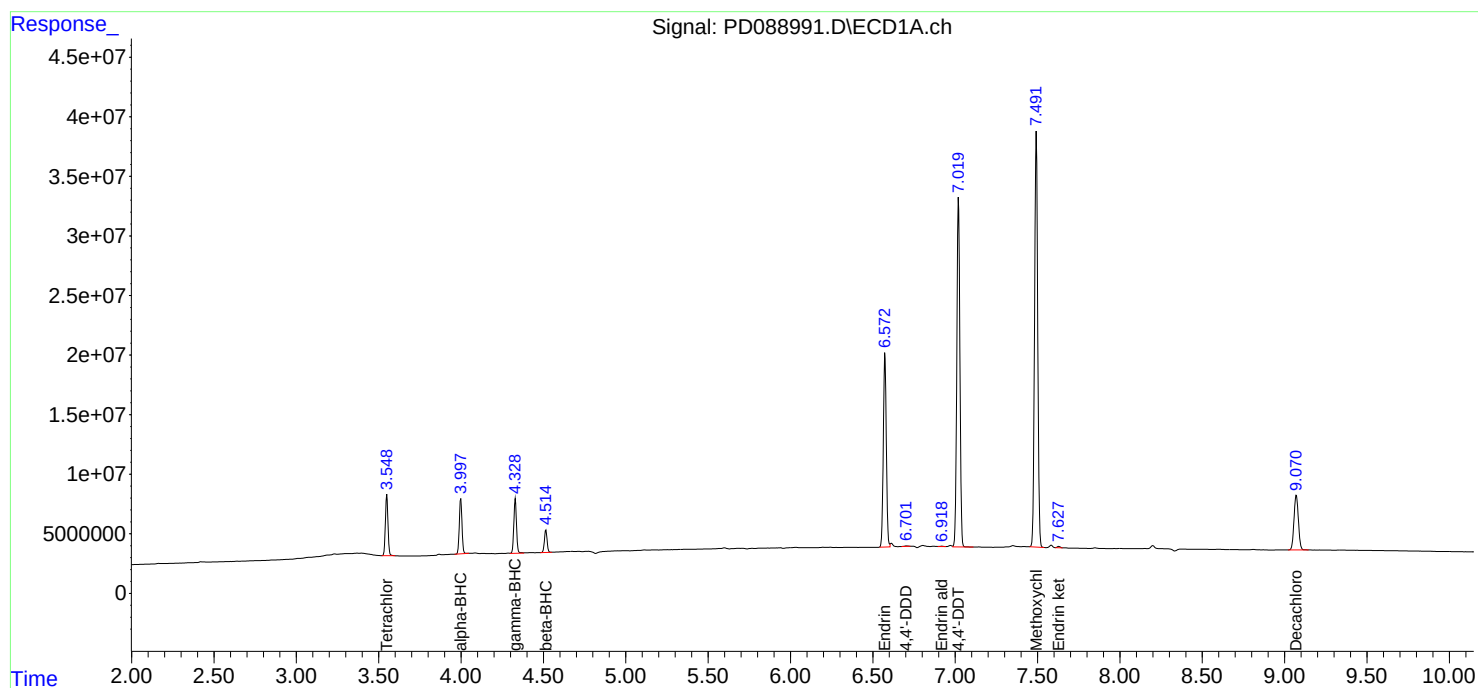
Instrument :
ECD_D
ClientSampleId :
PEM

Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 06/18/2025
Supervised By : mohammad ahmed 06/19/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:45:29 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089265.D Date Analyzed: 07/01/2025

Lab Sample No.(PEM): PEM Time Analyzed: 12:13

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.074	8.970	9.170	19.270	20.000	-3.7
Tetrachloro-m-xylene	3.550	3.500	3.600	18.530	20.000	-7.4
alpha-BHC	4.000	3.950	4.050	8.560	10.000	-14.4
beta-BHC	4.516	4.470	4.570	9.530	10.000	-4.7
gamma-BHC (Lindane)	4.329	4.280	4.380	8.840	10.000	-11.6
Endrin	6.573	6.500	6.640	43.050	50.000	-13.9
4,4'-DDT	7.022	6.950	7.090	78.670	100.000	-21.3
Methoxychlor	7.493	7.420	7.560	176.900	250.000	-29.2

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089265.D Date Analyzed: 07/01/2025

Lab Sample No.(PEM): PEM Time Analyzed: 12:13

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.072	7.970	8.170	19.650	20.000	-1.8
Tetrachloro-m-xylene	2.881	2.830	2.930	19.970	20.000	-0.2
alpha-BHC	3.392	3.340	3.440	10.130	10.000	1.3
beta-BHC	4.025	3.970	4.080	9.890	10.000	-1.1
gamma-BHC (Lindane)	3.730	3.680	3.780	9.870	10.000	-1.3
Endrin	5.789	5.720	5.860	45.160	50.000	-9.7
4,4'-DDT	6.183	6.110	6.250	78.010	100.000	-22.0
Methoxychlor	6.754	6.680	6.820	158.950	250.000	-36.4

PEM
 Data File: PD089265.D Date Acquired 7/1/2025 12:13
 Operator: AR\AJ

ENDRIN BREAK DOWN

Column #1

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin	6.57	177602871.5	188682387.7	11079516.2	Down 5.87
Endrin aldehyde	6.92	3048057.244			
Endrin ketone	7.63	8031458.925			

Column #2

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin #2	5.79	980555114.1	1082275727	101720613	9.40
Endrin aldehyde #2	6.26	27499938.02			
Endrin ketone #2	6.99	74220675.1			

DDT BREAK DOWN

Column #1

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT	7.02	298004872.3	317256503	19251630.6	6.07
4,4'-DDE	6.19	938270.216			
4,4'-DDD	6.70	18313360.43			

Column #2

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT #2	6.18	1581914044	1693486508	111572464	6.59
4,4'-DDE #2	5.37	5646016.685			
4,4'-DDD #2	5.93	105926447.1			

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089265.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 12:13
 Operator : AR\AJ
 Sample : PEM
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PEM

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:44:49 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	52857529	338.6E6	18.526	19.975
28)	SA Decachlor...	9.074	8.072	75689745	388.5E6	19.271	19.651
Target Compounds							
2)	A alpha-BHC	4.000	3.392	49881725	271.5E6	8.557	10.133m
3)	MA gamma-BHC...	4.329	3.730	49349393	244.4E6	8.841m	9.871
6)	B beta-BHC	4.516	4.025	21028261	106.2E6	9.527	9.887
12)	B 4,4'-DDE	6.194	5.374	938270	5646017	0.215m	0.246m
14)	MA Endrin	6.573	5.789	177.6E6	980.6E6	43.049m	45.158
16)	A 4,4'-DDD	6.704	5.929	18313360	105.9E6	5.336m	5.537
17)	MA 4,4'-DDT	7.022	6.183	298.0E6	1581.9E6	78.667	78.006
18)	B Endrin al...	6.915	6.258	3048057	27499938	0.995	1.796 #
20)	A Methoxychlor	7.493	6.754	357.5E6	1711.6E6	176.903	158.950
21)	B Endrin ke...	7.630	6.990	8031459	74220675	1.972	3.436 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089265.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 12:13
Operator : AR\AJ
Sample : PEM
Misc :
ALS Vial : 3 Sample Multiplier: 1

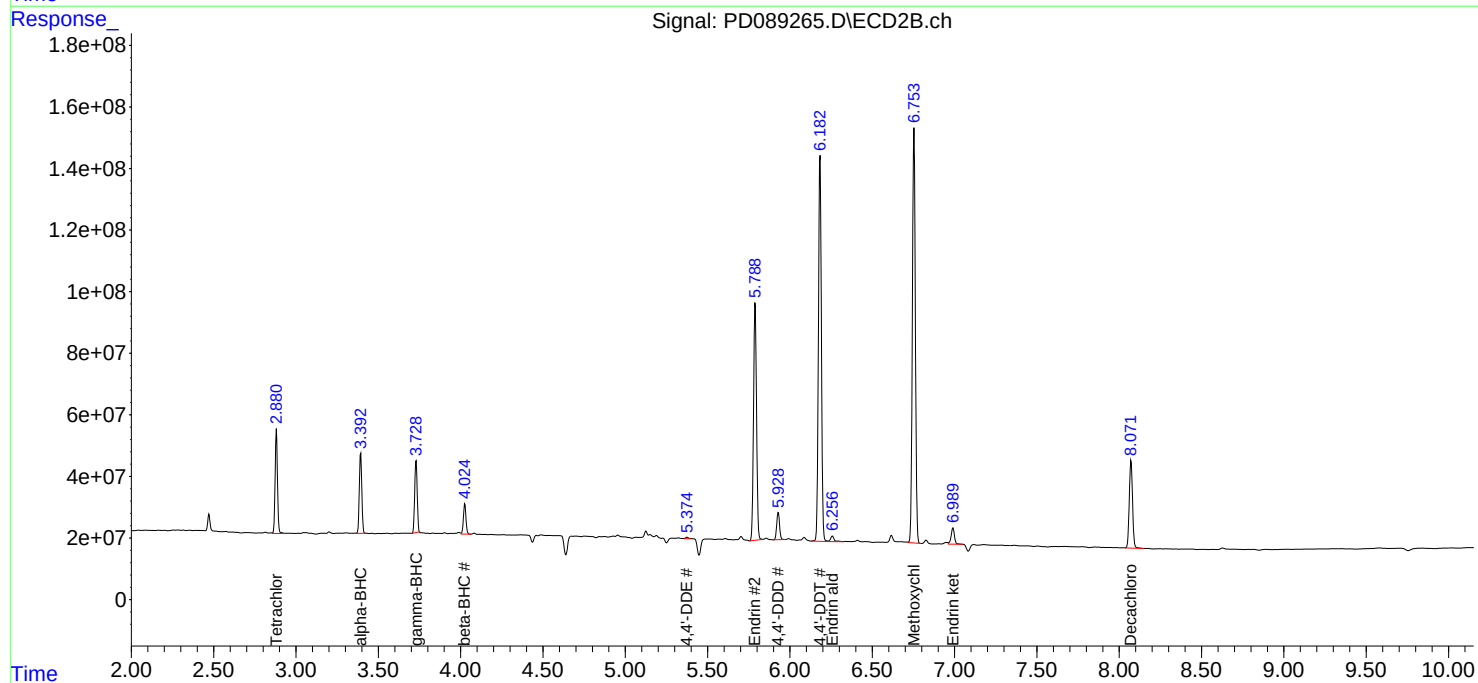
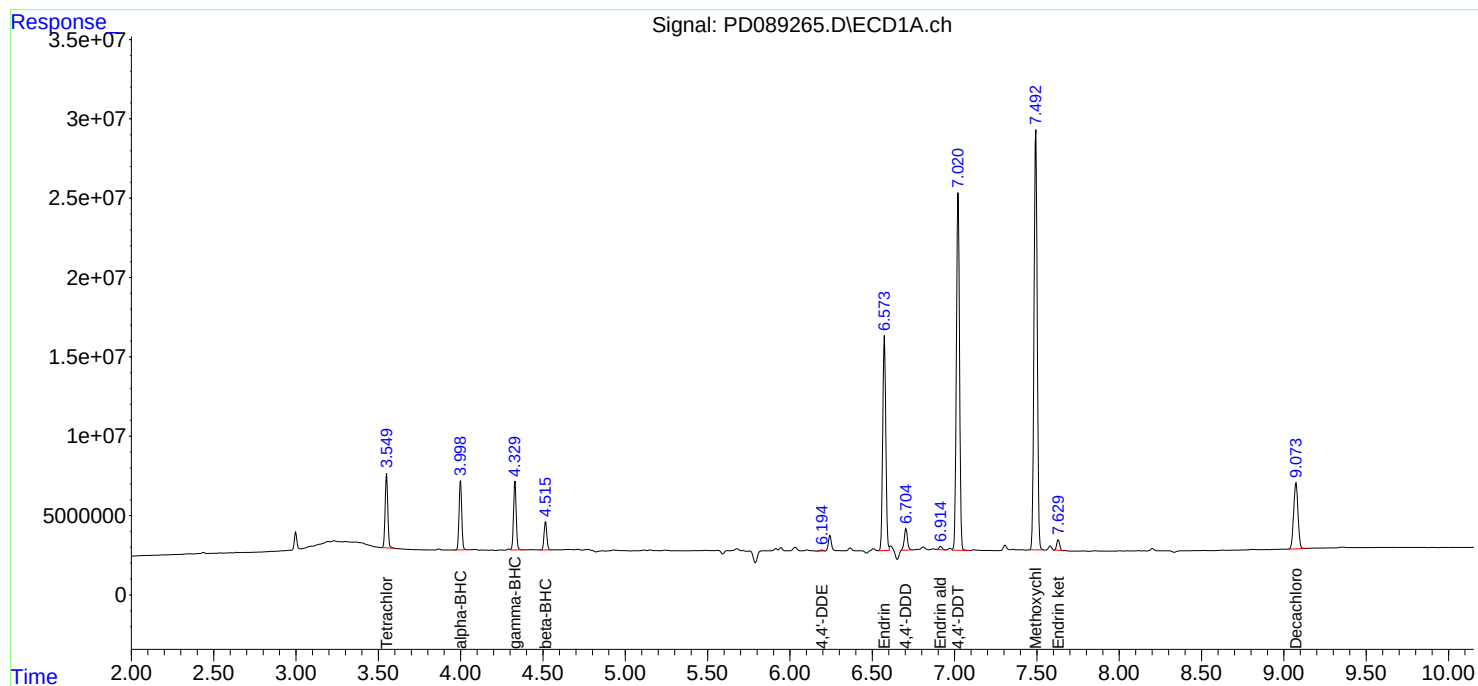
Instrument :
ECD_D
ClientSampleId :
PEM

Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:44:49 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089288.D Date Analyzed: 07/01/2025

Lab Sample No.(PEM): PEM Time Analyzed: 19:21

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.072	8.970	9.170	17.760	20.000	-11.2
Tetrachloro-m-xylene	3.550	3.500	3.600	18.810	20.000	-6.0
alpha-BHC	3.999	3.950	4.050	8.640	10.000	-13.6
beta-BHC	4.516	4.470	4.570	9.570	10.000	-4.3
gamma-BHC (Lindane)	4.330	4.280	4.380	9.110	10.000	-8.9
Endrin	6.572	6.500	6.640	42.570	50.000	-14.9
4,4'-DDT	7.021	6.950	7.090	74.530	100.000	-25.5
Methoxychlor	7.493	7.420	7.560	168.790	250.000	-32.5

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089288.D Date Analyzed: 07/01/2025

Lab Sample No.(PEM): PEM Time Analyzed: 19:21

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.071	7.970	8.170	17.280	20.000	-13.6
Tetrachloro-m-xylene	2.881	2.830	2.930	20.200	20.000	1.0
alpha-BHC	3.393	3.340	3.440	10.600	10.000	6.0
beta-BHC	4.025	3.970	4.080	9.990	10.000	-0.1
gamma-BHC (Lindane)	3.728	3.680	3.780	10.260	10.000	2.6
Endrin	5.789	5.720	5.860	44.950	50.000	-10.1
4,4'-DDT	6.183	6.110	6.250	73.250	100.000	-26.8
Methoxychlor	6.753	6.680	6.820	152.400	250.000	-39.0

Data File: PEM
 PD089288.D Date Acquired 7/1/2025 19:21
 Operator: AR\AJ

ENDRIN BREAK DOWN

Column #1

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin	6.57	175633744.9	185964995.6	10331250.7	Down 5.56
Endrin aldehyde	6.91	2694696.517			
Endrin ketone	7.63	7636554.187			

Column #2

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin #2	5.79	976002765.7	1056010998	80008232.4	7.58
Endrin aldehyde #2	6.26	18199292.53			
Endrin ketone #2	6.99	61808939.84			

DDT BREAK DOWN

Column #1

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT	7.02	282343090.9	305571082.9	23227992	7.60
4,4'-DDE	6.19	659131.264			
4,4'-DDD	6.70	22568860.72			

Column #2

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT #2	6.18	1485498279	1602605944	117107665	7.31
4,4'-DDE #2	5.37	5365272.746			
4,4'-DDD #2	5.93	111742391.8			

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089288.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 19:21
 Operator : AR\AJ
 Sample : PEM
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PEM

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:50:34 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	53669144	342.4E6	18.811	20.196
28)	SA Decachlor...	9.072	8.071	69751872	341.6E6	17.759	17.278
Target Compounds							
2)	A alpha-BHC	3.999	3.393	50380963	283.9E6	8.643	10.595
3)	MA gamma-BHC...	4.330	3.728	50837015	254.1E6	9.108	10.263m
6)	B beta-BHC	4.516	4.025	21114096	107.4E6	9.566	9.993
12)	B 4,4'-DDE	6.193	5.374	659131	5365273	0.151	0.234m#
14)	MA Endrin	6.572	5.789	175.6E6	976.0E6	42.572m	44.948
16)	A 4,4'-DDD	6.704	5.929	22568861	111.7E6	6.576	5.841
17)	MA 4,4'-DDT	7.021	6.183	282.3E6	1485.5E6	74.532	73.252
18)	B Endrin al...	6.915	6.257	2694697	18199293	0.879	1.189 #
20)	A Methoxychlor	7.493	6.753	341.1E6	1641.0E6	168.786	152.396
21)	B Endrin ke...	7.629	6.990	7636554	61808940	1.875	2.861 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089288.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 19:21
Operator : AR\AJ
Sample : PEM
Misc :
ALS Vial : 3 Sample Multiplier: 1

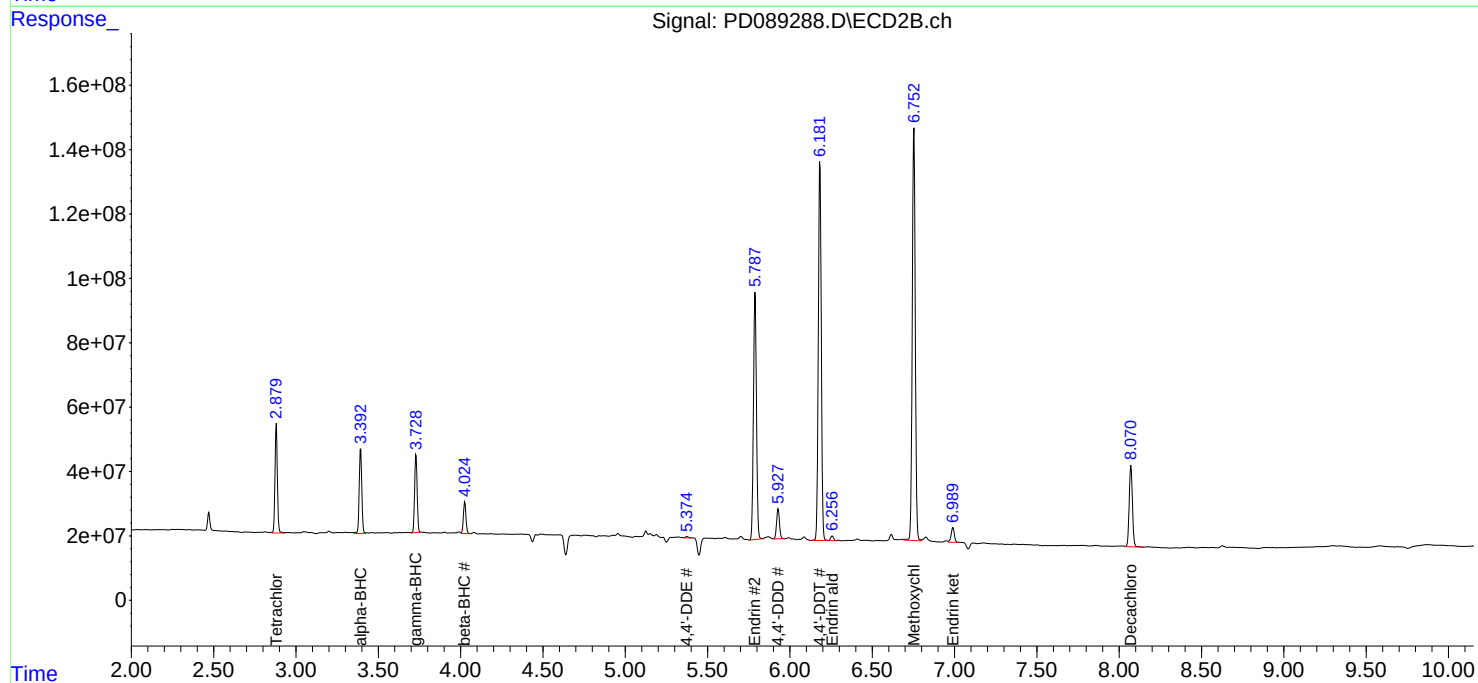
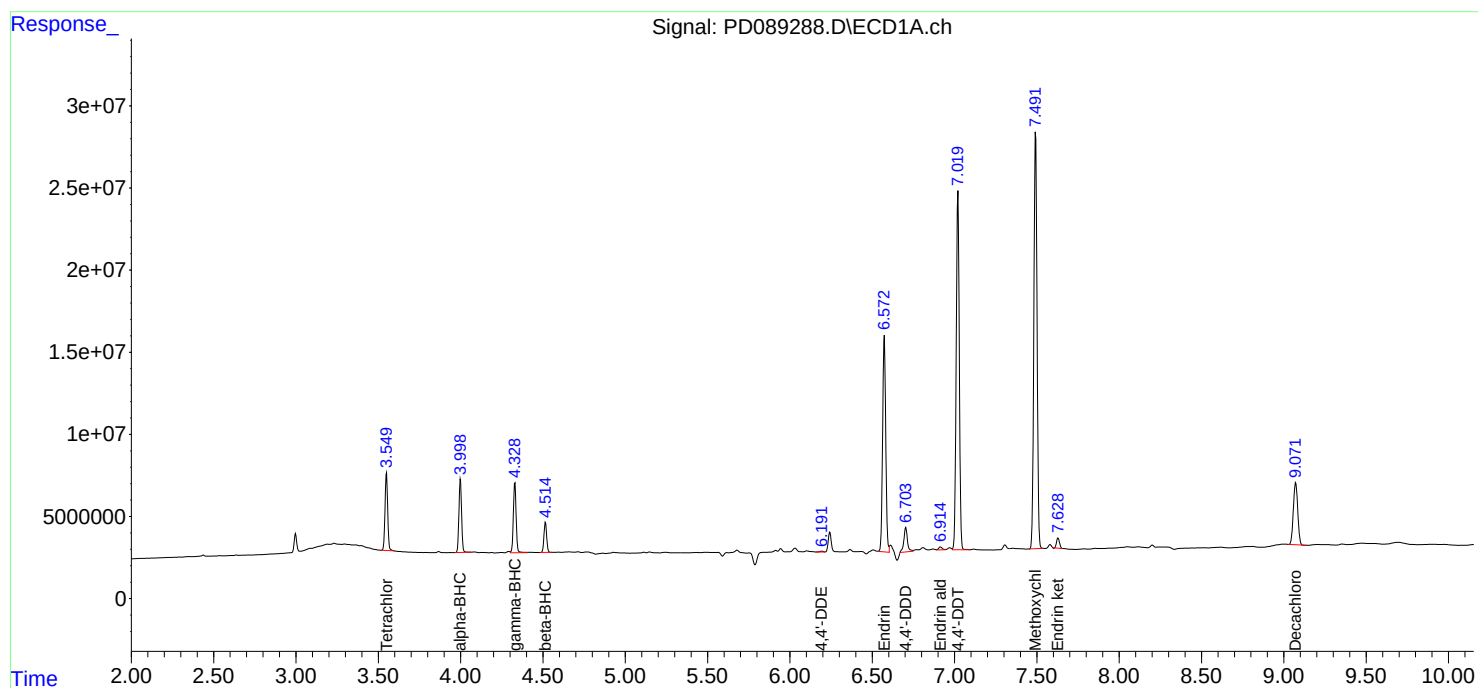
Instrument :
ECD_D
ClientSampleId :
PEM

Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/02/2025
Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:50:34 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089327.D Date Analyzed: 07/03/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:32

PEM COMPOUND	RT	RT WINDOW		CALC	NOM	%D
		FROM	TO	AMOUNT(ng)	AMOUNT(ng)	
Decachlorobiphenyl	9.073	8.970	9.170	19.550	20.000	-2.3
Tetrachloro-m-xylene	3.551	3.500	3.600	21.620	20.000	8.1
alpha-BHC	3.999	3.950	4.050	10.140	10.000	1.4
beta-BHC	4.516	4.470	4.570	11.040	10.000	10.4
gamma-BHC (Lindane)	4.331	4.280	4.380	10.550	10.000	5.5
Endrin	6.574	6.500	6.640	48.050	50.000	-3.9
4,4'-DDT	7.021	6.950	7.090	83.070	100.000	-16.9
Methoxychlor	7.494	7.420	7.560	183.850	250.000	-26.5

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 06/17/2025 06/17/2025

Client Sample No. (PEM): PEM - PD089327.D Date Analyzed: 07/03/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:32

PEM COMPOUND	RT	RT WINDOW		CALC	NOM	%D
		FROM	TO	AMOUNT(ng)	AMOUNT(ng)	
Decachlorobiphenyl	8.071	7.970	8.170	19.330	20.000	-3.4
Tetrachloro-m-xylene	2.881	2.830	2.930	23.640	20.000	18.2
alpha-BHC	3.393	3.340	3.440	12.290	10.000	22.9
beta-BHC	4.025	3.970	4.080	11.540	10.000	15.4
gamma-BHC (Lindane)	3.729	3.680	3.780	11.780	10.000	17.8
Endrin	5.788	5.720	5.860	51.080	50.000	2.2
4,4'-DDT	6.182	6.110	6.250	82.740	100.000	-17.3
Methoxychlor	6.753	6.680	6.820	160.450	250.000	-35.8

Data File: PEM
 PD089327.D Date Acquired 7/3/2025 9:32
 Operator: AR\AJ

ENDRIN BREAK DOWN

Column #1

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin	6.57	198227503.4	211531157.7	13303654.3	Down 6.29
Endrin aldehyde	6.92	3278521.382			
Endrin ketone	7.63	10025132.94			

Column #2

Name	RT	Response	Response [E+EA+EK]	Response [EA+EK]	% Break Down
Endrin #2	5.79	1109184579	1214927492	105742913	8.70
Endrin aldehyde #2	6.26	24041555.25			
Endrin ketone #2	6.99	81701357.78			

DDT BREAK DOWN

Column #1

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT	7.02	314699780.6	339367790.3	24668009.7	7.27
4,4'-DDE	6.19	899767.037			
4,4'-DDD	6.70	23768242.68			

Column #2

Name	RT	Response	Response [DDT+DDE+DDD]	Response [DDE+DDD]	% Break Down
4,4'-DDT #2	6.18	1677864534	1845597601	167733067	9.09
4,4'-DDE #2	5.37	7625893.265			
4,4'-DDD #2	5.93	160107173.2			

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089327.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 09:32
 Operator : AR\AJ
 Sample : PEM
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PEM

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/07/2025
 Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:16:52 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.881	61680202	400.7E6	21.618	23.636
28)	SA Decachlor...	9.073	8.071	76793160	382.1E6	19.552	19.329
Target Compounds							
2)	A alpha-BHC	3.999	3.393	59112351	329.4E6	10.141	12.293
3)	MA gamma-BHC...	4.331	3.729	58900726	291.7E6	10.553	11.781
6)	B beta-BHC	4.516	4.025	24368122	124.0E6	11.041	11.536
12)	B 4,4'-DDE	6.192	5.372	899767	7625893	0.206m	0.333m#
14)	MA Endrin	6.574	5.788	198.2E6	1109.2E6	48.048	51.082
16)	A 4,4'-DDD	6.704	5.927	23768243	160.1E6	6.925m	8.370m
17)	MA 4,4'-DDT	7.021	6.182	314.7E6	1677.9E6	83.074	82.738
18)	B Endrin al...	6.916	6.257	3278521	24041555	1.070	1.570 #
20)	A Methoxychlor	7.494	6.753	371.5E6	1727.7E6	183.849	160.450
21)	B Endrin ke...	7.630	6.989	10025133	81701358	2.461	3.782 #

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089327.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 09:32
Operator : AR\AJ
Sample : PEM
Misc :
ALS Vial : 3 Sample Multiplier: 1

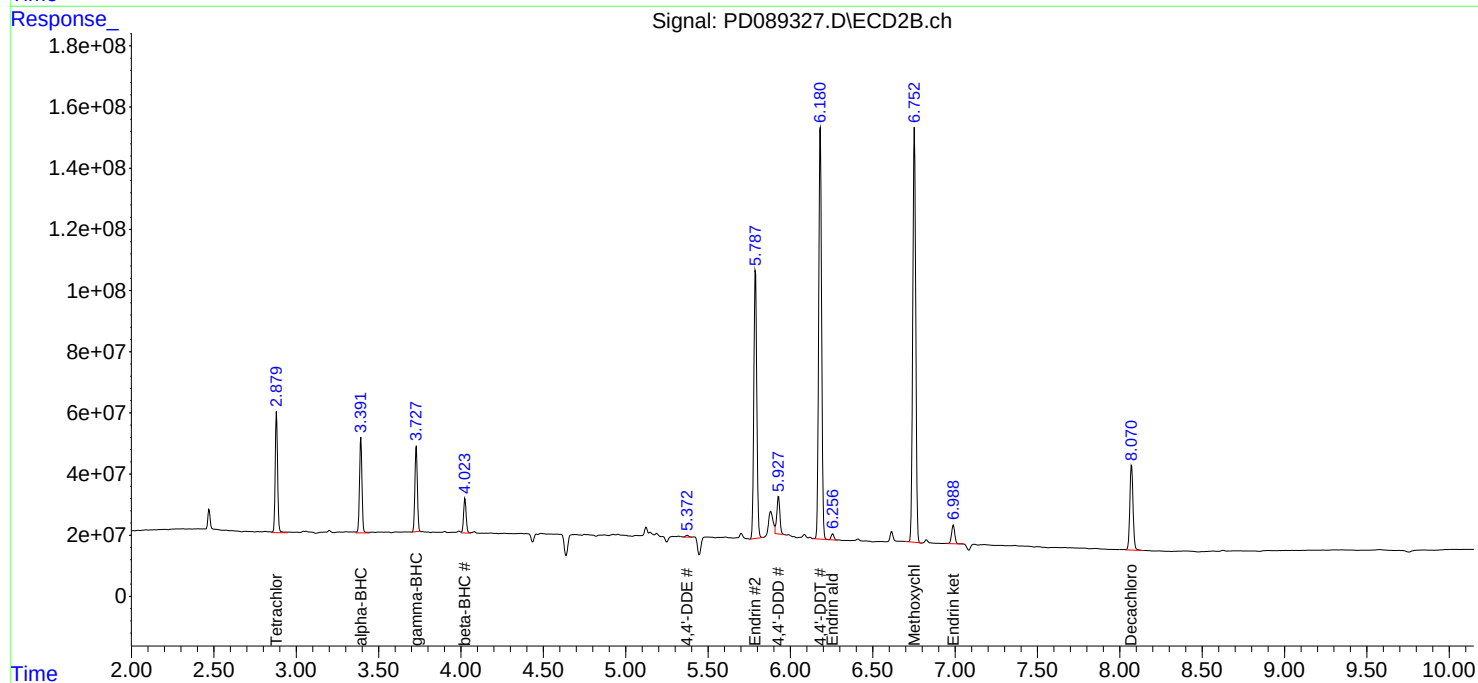
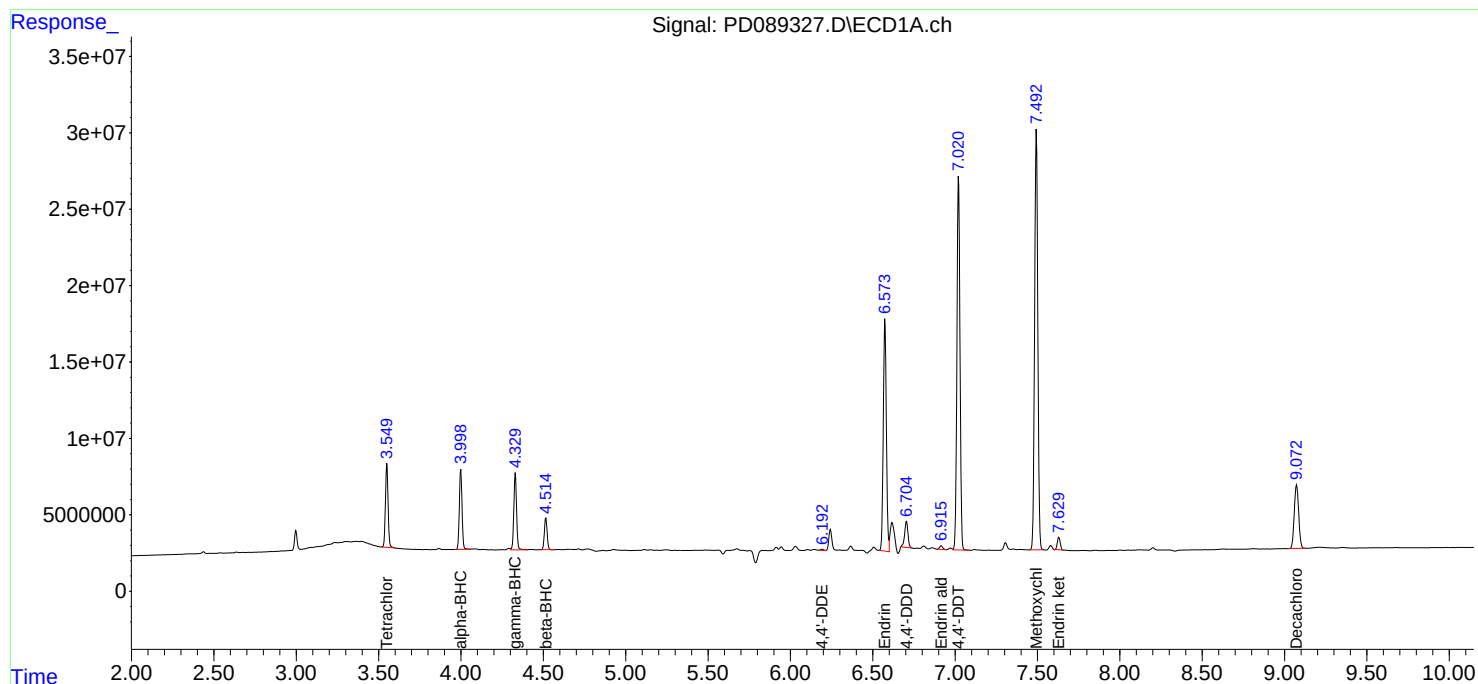
Instrument :
ECD_D
ClientSampleId :
PEM

Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/07/2025
Supervised By : mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:16:52 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Analytical Sequence

Client: CDM Smith	SDG No.: Q2458
Project: South River WM Replacement	Instrument ID: ECD_D
GC Column: ZB-MR1	ID: 0.32 (mm) Inst. Calib. Date(s): 06/17/2025 06/17/2025

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES,
AND STANDARDS IS GIVEN BELOW:

CLIENT ID	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	DATAFILE	DCB RT #	TCX RT #
LBLK	LBLK	06/17/2025	15:11	PD088990.D	9.07	3.55
PEM	PEM	06/17/2025	15:25	PD088991.D	9.07	3.55
RESCHK	RESCHK	06/17/2025	15:39	PD088992.D	9.07	3.55
PSTDICCC100	PSTDICCC100	06/17/2025	15:52	PD088993.D	9.07	3.55
PSTDICCC075	PSTDICCC075	06/17/2025	16:06	PD088994.D	9.07	3.55
PSTDICCC050	PSTDICCC050	06/17/2025	16:20	PD088995.D	9.07	3.55
PSTDICCC025	PSTDICCC025	06/17/2025	16:33	PD088996.D	9.07	3.55
PSTDICCC005	PSTDICCC005	06/17/2025	16:47	PD088997.D	9.07	3.55
PCHLORICC500	PCHLORICC500	06/17/2025	17:28	PD089000.D	9.07	3.55
PTOXICC500	PTOXICC500	06/17/2025	18:36	PD089005.D	9.07	3.55
LBLK	LBLK	07/01/2025	11:59	PD089264.D	9.07	3.55
PEM	PEM	07/01/2025	12:13	PD089265.D	9.07	3.55
PSTDCCC050	PSTDCCC050	07/01/2025	13:35	PD089266.D	9.08	3.56
PB168672BL	PB168672BL	07/01/2025	14:33	PD089270.D	9.07	3.55
PB168672BS	PB168672BS	07/01/2025	16:23	PD089275.D	9.08	3.56
LBLK	LBLK	07/01/2025	16:37	PD089276.D	9.07	3.55
PSTDCCC050	PSTDCCC050	07/01/2025	16:50	PD089277.D	9.07	3.55
TP-76	Q2458-01	07/01/2025	18:26	PD089284.D	9.07	3.55
TP-55	Q2458-02	07/01/2025	18:40	PD089285.D	9.07	3.55
TP-68	Q2458-03	07/01/2025	18:53	PD089286.D	9.07	3.55
LBLK	LBLK	07/01/2025	19:07	PD089287.D	9.07	3.55
PEM	PEM	07/01/2025	19:21	PD089288.D	9.07	3.55
PSTDCCC050	PSTDCCC050	07/01/2025	19:34	PD089289.D	9.07	3.55
TP-67	Q2458-04	07/01/2025	20:15	PD089290.D	9.07	3.55
TP-67MS	Q2458-04MS	07/01/2025	20:29	PD089291.D	9.07	3.55
TP-67MSD	Q2458-04MSD	07/01/2025	20:43	PD089292.D	9.07	3.55
TP-66	Q2458-05	07/01/2025	20:57	PD089293.D	9.07	3.55
TP-60	Q2458-06	07/01/2025	21:10	PD089294.D	9.07	3.55
TP-62	Q2458-07	07/01/2025	21:24	PD089295.D	9.07	3.55
TP-63	Q2458-08	07/01/2025	21:37	PD089296.D	9.07	3.55
TP-59	Q2458-09	07/01/2025	21:51	PD089297.D	9.07	3.55
LBLK	LBLK	07/01/2025	22:05	PD089298.D	9.07	3.55
PSTDCCC050	PSTDCCC050	07/01/2025	22:59	PD089299.D	9.07	3.55
LBLK	LBLK	07/03/2025	09:19	PD089326.D	9.08	3.55
PEM	PEM	07/03/2025	09:32	PD089327.D	9.07	3.55
PSTDCCC050	PSTDCCC050	07/03/2025	09:46	PD089328.D	9.07	3.55
PB168718BL	PB168718BL	07/03/2025	13:55	PD089329.D	9.08	3.56
PB168718BS	PB168718BS	07/03/2025	14:09	PD089330.D	9.08	3.55
PB168718BSD	PB168718BSD	07/03/2025	14:27	PD089331.D	9.08	3.56
FB-06272025	Q2458-10	07/03/2025	14:40	PD089332.D	9.08	3.55
LBLK	LBLK	07/03/2025	15:48	PD089337.D	9.08	3.55
PSTDCCC050	PSTDCCC050	07/03/2025	16:02	PD089338.D	9.07	3.55

Analytical Sequence

Analytical Sequence

Client: CDM Smith

SDG No.: Q2458

Project: South River WM Replacement

Instrument ID: ECD_D

GC Column: ZB-MR2

ID: 0.32 (mm)

Inst. Calib. Date(s): 06/17/2025

06/17/2025

THE ANALYTICAL SEQUENCE OF PERFORMANCE EVALUATION MIXTURES, BLANKS, SAMPLES,
AND STANDARDS IS GIVEN BELOW:

CLIENT ID	LAB SAMPLE ID	DATE ANALYZED	TIME ANALYZED	DATAFILE	DCB RT #	TCX RT #
IBLK	IBLK	06/17/2025	15:11	PD088990.D	8.07	2.88
PEM	PEM	06/17/2025	15:25	PD088991.D	8.07	2.88
RESCHK	RESCHK	06/17/2025	15:39	PD088992.D	8.07	2.88
PSTDICCC100	PSTDICCC100	06/17/2025	15:52	PD088993.D	8.07	2.88
PSTDICCC075	PSTDICCC075	06/17/2025	16:06	PD088994.D	8.07	2.88
PSTDICCC050	PSTDICCC050	06/17/2025	16:20	PD088995.D	8.07	2.88
PSTDICCC025	PSTDICCC025	06/17/2025	16:33	PD088996.D	8.07	2.88
PSTDICCC005	PSTDICCC005	06/17/2025	16:47	PD088997.D	8.07	2.88
PCHLORICCC500	PCHLORICCC500	06/17/2025	17:28	PD089000.D	8.07	2.88
PTOXICCC500	PTOXICCC500	06/17/2025	18:36	PD089005.D	8.07	2.88
IBLK	IBLK	07/01/2025	11:59	PD089264.D	8.07	2.88
PEM	PEM	07/01/2025	12:13	PD089265.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/01/2025	13:35	PD089266.D	8.07	2.88
PB168672BL	PB168672BL	07/01/2025	14:33	PD089270.D	8.07	2.88
PB168672BS	PB168672BS	07/01/2025	16:23	PD089275.D	8.07	2.88
IBLK	IBLK	07/01/2025	16:37	PD089276.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/01/2025	16:50	PD089277.D	8.07	2.88
TP-76	Q2458-01	07/01/2025	18:26	PD089284.D	8.07	2.88
TP-55	Q2458-02	07/01/2025	18:40	PD089285.D	8.07	2.88
TP-68	Q2458-03	07/01/2025	18:53	PD089286.D	8.07	2.88
IBLK	IBLK	07/01/2025	19:07	PD089287.D	8.07	2.88
PEM	PEM	07/01/2025	19:21	PD089288.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/01/2025	19:34	PD089289.D	8.07	2.88
TP-67	Q2458-04	07/01/2025	20:15	PD089290.D	8.07	2.88
TP-67MS	Q2458-04MS	07/01/2025	20:29	PD089291.D	8.07	2.88
TP-67MSD	Q2458-04MSD	07/01/2025	20:43	PD089292.D	8.07	2.88
TP-66	Q2458-05	07/01/2025	20:57	PD089293.D	8.07	2.88
TP-60	Q2458-06	07/01/2025	21:10	PD089294.D	8.07	2.88
TP-62	Q2458-07	07/01/2025	21:24	PD089295.D	8.07	2.88
TP-63	Q2458-08	07/01/2025	21:37	PD089296.D	8.07	2.88
TP-59	Q2458-09	07/01/2025	21:51	PD089297.D	8.07	2.88
IBLK	IBLK	07/01/2025	22:05	PD089298.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/01/2025	22:59	PD089299.D	8.07	2.88
IBLK	IBLK	07/03/2025	09:19	PD089326.D	8.07	2.88
PEM	PEM	07/03/2025	09:32	PD089327.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/03/2025	09:46	PD089328.D	8.07	2.88
PB168718BL	PB168718BL	07/03/2025	13:55	PD089329.D	8.08	2.88
PB168718BS	PB168718BS	07/03/2025	14:09	PD089330.D	8.07	2.88
PB168718BSD	PB168718BSD	07/03/2025	14:27	PD089331.D	8.07	2.88
FB-06272025	Q2458-10	07/03/2025	14:40	PD089332.D	8.07	2.88
IBLK	IBLK	07/03/2025	15:48	PD089337.D	8.07	2.88
PSTDCCC050	PSTDCCC050	07/03/2025	16:02	PD089338.D	8.07	2.88

Analytical Sequence

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168672BS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168672BS

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDD	1	6.71	6.66	6.76	18.7	4.4
	2	5.93	5.88	5.98	17.9	
4,4'-DDT	1	7.03	6.98	7.08	16.3	2.5
	2	6.19	6.14	6.24	15.9	
alpha-BHC	1	4.01	3.96	4.06	18.7	0.5
	2	3.39	3.34	3.44	18.6	
Aldrin	1	5.28	5.23	5.33	18.7	1.6
	2	4.37	4.32	4.42	18.4	
beta-BHC	1	4.52	4.47	4.57	17.7	1.7
	2	4.03	3.98	4.08	18.0	
alpha-Chlordane	1	6.03	5.98	6.08	18.1	0.6
	2	5.19	5.14	5.24	18.2	
4,4'-DDE	1	6.20	6.15	6.25	18.4	2.8
	2	5.38	5.33	5.43	17.9	
Endosulfan II	1	6.79	6.74	6.84	17.9	0.6
	2	6.08	6.03	6.13	17.8	
Endrin aldehyde	1	6.92	6.87	6.97	17.8	1.1
	2	6.26	6.21	6.31	17.6	
Endosulfan sulfate	1	7.16	7.11	7.21	17.7	0.6
	2	6.48	6.43	6.53	17.6	
Methoxychlor	1	7.50	7.45	7.55	16.3	1.8
	2	6.76	6.71	6.81	16.6	
Endrin ketone	1	7.64	7.59	7.69	18.3	2.8
	2	6.99	6.94	7.04	17.8	
gamma-BHC (Lindane)	1	4.34	4.29	4.39	18.6	1.1
	2	3.73	3.68	3.78	18.4	
Heptachlor	1	4.94	4.89	4.99	18.2	3.4
	2	4.08	4.03	4.13	17.6	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168672BS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168672BS

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
delta-BHC	1	4.77	4.72	4.82	19.2	4.8
	2	4.26	4.21	4.31	18.3	
Heptachlor epoxide	1	5.70	5.65	5.75	18.0	7
	2	4.87	4.82	4.92	19.3	
Endosulfan I	1	6.08	6.03	6.13	18.3	0.5
	2	5.25	5.20	5.30	18.4	
gamma-Chlordane	1	5.95	5.90	6.00	18.5	3.3
	2	5.13	5.08	5.18	17.9	
Dieldrin	1	6.35	6.30	6.40	18.4	2.2
	2	5.51	5.46	5.56	18.0	
Endrin	1	6.58	6.53	6.63	16.5	0
	2	5.79	5.74	5.84	16.5	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168718BS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168718BS

Date(s) Analyzed: 07/03/2025 07/03/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column: (2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
Endrin aldehyde	1	6.92	6.87	6.97	0.56	1.7
	2	6.26	6.21	6.31	0.55	
Methoxychlor	1	7.50	7.45	7.55	0.42	5.3
	2	6.75	6.70	6.80	0.44	
Endrin ketone	1	7.63	7.58	7.68	0.56	0.6
	2	6.99	6.94	7.04	0.55	
gamma-BHC (Lindane)	1	4.33	4.28	4.38	0.55	1
	2	3.73	3.68	3.78	0.56	
Heptachlor	1	4.93	4.88	4.98	0.54	2.3
	2	4.08	4.03	4.13	0.52	
Heptachlor epoxide	1	5.69	5.64	5.74	0.55	0.5
	2	4.87	4.82	4.92	0.55	
gamma-Chlordane	1	5.95	5.90	6.00	0.57	4.4
	2	5.13	5.08	5.18	0.55	
Endrin	1	6.58	6.53	6.63	0.49	2.8
	2	5.79	5.74	5.84	0.48	
4,4'-DDD	1	6.71	6.66	6.76	0.59	5.1
	2	5.93	5.88	5.98	0.56	
4,4'-DDE	1	6.20	6.15	6.25	0.56	2.5
	2	5.38	5.33	5.43	0.55	
4,4'-DDT	1	7.02	6.97	7.07	0.48	1.3
	2	6.18	6.13	6.23	0.47	
alpha-BHC	1	4.00	3.95	4.05	0.56	1
	2	3.39	3.34	3.44	0.56	
Aldrin	1	5.27	5.22	5.32	0.56	1.3
	2	4.37	4.32	4.42	0.56	
alpha-Chlordane	1	6.03	5.98	6.08	0.56	1.9
	2	5.19	5.14	5.24	0.55	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168718BS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168718BS

Date(s) Analyzed: 07/03/2025 07/03/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
Endosulfan II	1	6.79	6.74	6.84	0.57	3.2
	2	6.08	6.03	6.13	0.55	
Endosulfan sulfate	1	7.15	7.10	7.20	0.55	0.6
	2	6.48	6.43	6.53	0.55	
beta-BHC	1	4.52	4.47	4.57	0.53	2.8
	2	4.03	3.98	4.08	0.54	
delta-BHC	1	4.77	4.72	4.82	0.59	6
	2	4.26	4.21	4.31	0.55	
Endosulfan I	1	6.08	6.03	6.13	0.56	0.8
	2	5.25	5.20	5.30	0.55	
Dieldrin	1	6.35	6.30	6.40	0.57	2.4
	2	5.51	5.46	5.56	0.55	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168718BSD

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168718BSD

Date(s) Analyzed: 07/03/2025 07/03/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDD	1	6.71	6.66	6.76	0.58	4.1
	2	5.93	5.88	5.98	0.56	
4,4'-DDT	1	7.03	6.98	7.08	0.47	0.6
	2	6.18	6.13	6.23	0.47	
alpha-BHC	1	4.01	3.96	4.06	0.55	1.1
	2	3.39	3.34	3.44	0.56	
Aldrin	1	5.28	5.23	5.33	0.56	0.6
	2	4.37	4.32	4.42	0.55	
beta-BHC	1	4.52	4.47	4.57	0.52	2.8
	2	4.03	3.98	4.08	0.54	
alpha-Chlordane	1	6.03	5.98	6.08	0.55	1.4
	2	5.19	5.14	5.24	0.54	
4,4'-DDE	1	6.20	6.15	6.25	0.55	1.2
	2	5.38	5.33	5.43	0.55	
Endosulfan II	1	6.79	6.74	6.84	0.56	2.7
	2	6.08	6.03	6.13	0.54	
Endrin aldehyde	1	6.92	6.87	6.97	0.55	0.8
	2	6.26	6.21	6.31	0.55	
Endosulfan sulfate	1	7.16	7.11	7.21	0.55	0.1
	2	6.48	6.43	6.53	0.55	
Methoxychlor	1	7.50	7.45	7.55	0.42	5.6
	2	6.76	6.71	6.81	0.44	
Endrin ketone	1	7.64	7.59	7.69	0.55	0.6
	2	6.99	6.94	7.04	0.55	
gamma-BHC (Lindane)	1	4.34	4.29	4.39	0.55	1.2
	2	3.73	3.68	3.78	0.55	
Heptachlor	1	4.94	4.89	4.99	0.53	1.4
	2	4.08	4.03	4.13	0.52	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB168718BSD

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: PB168718BSD

Date(s) Analyzed: 07/03/2025 07/03/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
delta-BHC	1	4.77	4.72	4.82	0.57	3.9
	2	4.26	4.21	4.31	0.55	
Heptachlor epoxide	1	5.70	5.65	5.75	0.54	1.5
	2	4.87	4.82	4.92	0.55	
Endosulfan I	1	6.08	6.03	6.13	0.55	0.1
	2	5.25	5.20	5.30	0.55	
gamma-Chlordane	1	5.95	5.90	6.00	0.56	3.3
	2	5.13	5.08	5.18	0.55	
Dieldrin	1	6.35	6.30	6.40	0.56	2
	2	5.51	5.46	5.56	0.55	
Endrin	1	6.58	6.53	6.63	0.48	1
	2	5.79	5.74	5.84	0.48	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-60

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-06

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDT	1	7.02	6.97	7.07	0.15	80.9
	2	6.18	6.13	6.23	0.36	
gamma-Chlordane	1	5.94	5.89	5.99	0.46	35.9
	2	5.12	5.07	5.17	0.32	
alpha-Chlordane	1	6.03	5.98	6.08	0.90	43.7
	2	5.19	5.14	5.24	0.58	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-63

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-08

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDD	1	6.70	6.65	6.75	0.41	25.2
	2	5.93	5.88	5.98	0.32	
4,4'-DDT	1	7.02	6.97	7.07	0.24	57.6
	2	6.18	6.13	6.23	0.44	
gamma-Chlordane	1	5.94	5.89	5.99	0.40	31.4
	2	5.12	5.07	5.17	0.29	
alpha-Chlordane	1	6.03	5.98	6.08	0.58	22.7
	2	5.19	5.14	5.24	0.47	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-66

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-05

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDT	1	7.02	6.97	7.07	0.24	48.8
	2	6.18	6.13	6.23	0.40	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-67

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-04

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
gamma-Chlordane	1	5.94	5.89	5.99	0.29	20
	2	5.13	5.08	5.18	0.23	
alpha-Chlordane	1	6.03	5.98	6.08	0.48	48.1
	2	5.19	5.14	5.24	0.29	
4,4'-DDE	1	6.20	6.15	6.25	0.21	49.7
	2	5.37	5.32	5.42	0.34	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-67MS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-04MS

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDD	1	6.71	6.66	6.76	18.4	7.9
	2	5.93	5.88	5.98	17.0	
4,4'-DDT	1	7.02	6.97	7.07	14.0	1.4
	2	6.18	6.13	6.23	14.2	
4,4'-DDE	1	6.20	6.15	6.25	18.0	0.6
	2	5.38	5.33	5.43	17.9	
Endosulfan II	1	6.79	6.74	6.84	17.1	0.6
	2	6.08	6.03	6.13	17.2	
Endrin aldehyde	1	6.92	6.87	6.97	16.8	1.2
	2	6.26	6.21	6.31	16.6	
Endosulfan sulfate	1	7.15	7.10	7.20	16.4	1.8
	2	6.48	6.43	6.53	16.1	
Methoxychlor	1	7.49	7.44	7.54	13.4	0.7
	2	6.75	6.70	6.80	13.3	
Endrin ketone	1	7.63	7.58	7.68	17.2	5.4
	2	6.99	6.94	7.04	16.3	
alpha-BHC	1	4.00	3.95	4.05	17.7	0
	2	3.39	3.34	3.44	17.7	
gamma-BHC (Lindane)	1	4.33	4.28	4.38	17.5	0
	2	3.73	3.68	3.78	17.5	
Heptachlor	1	4.93	4.88	4.98	16.5	1.2
	2	4.08	4.03	4.13	16.3	
Aldrin	1	5.27	5.22	5.32	18.0	0.6
	2	4.37	4.32	4.42	17.9	
beta-BHC	1	4.52	4.47	4.57	17.3	1.7
	2	4.03	3.98	4.08	17.6	
delta-BHC	1	4.76	4.71	4.81	17.9	5.7
	2	4.26	4.21	4.31	16.9	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-67MS

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-04MS

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
Heptachlor epoxide	1	5.69	5.64	5.74	17.6	3.4
	2	4.87	4.82	4.92	18.2	
Endosulfan I	1	6.07	6.02	6.12	17.8	1.1
	2	5.25	5.20	5.30	17.6	
gamma-Chlordane	1	5.94	5.89	5.99	18.2	0.5
	2	5.13	5.08	5.18	18.3	
alpha-Chlordane	1	6.03	5.98	6.08	18.0	0.6
	2	5.19	5.14	5.24	17.9	
Dieldrin	1	6.35	6.30	6.40	17.6	0.6
	2	5.51	5.46	5.56	17.7	
Endrin	1	6.57	6.52	6.62	16.6	1.8
	2	5.79	5.74	5.84	16.9	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-67MSD

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-04MSD

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column: (2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
Endosulfan II	1	6.79	6.74	6.84	17.3	0.6
	2	6.08	6.03	6.13	17.4	
4,4'-DDD	1	6.71	6.66	6.76	19.0	8.8
	2	5.93	5.88	5.98	17.4	
4,4'-DDT	1	7.02	6.97	7.07	14.7	2
	2	6.18	6.13	6.23	15.0	
Endrin aldehyde	1	6.92	6.87	6.97	17.3	0.6
	2	6.26	6.21	6.31	17.2	
Endosulfan sulfate	1	7.15	7.10	7.20	17.1	1.2
	2	6.48	6.43	6.53	16.9	
Methoxychlor	1	7.49	7.44	7.54	13.8	2.9
	2	6.75	6.70	6.80	14.2	
Endrin ketone	1	7.63	7.58	7.68	17.7	4
	2	6.99	6.94	7.04	17.0	
alpha-BHC	1	4.00	3.95	4.05	18.2	0.6
	2	3.39	3.34	3.44	18.1	
gamma-BHC (Lindane)	1	4.33	4.28	4.38	18.2	1.1
	2	3.73	3.68	3.78	18.0	
Heptachlor	1	4.93	4.88	4.98	16.9	0.6
	2	4.08	4.03	4.13	16.8	
Aldrin	1	5.27	5.22	5.32	18.2	0.6
	2	4.37	4.32	4.42	18.1	
beta-BHC	1	4.52	4.47	4.57	17.4	2.8
	2	4.03	3.98	4.08	17.9	
delta-BHC	1	4.76	4.71	4.81	19.0	6.5
	2	4.26	4.21	4.31	17.8	
Heptachlor epoxide	1	5.69	5.64	5.74	17.7	3.3
	2	4.87	4.82	4.92	18.3	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-67MSD

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-04MSD

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column: (2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
Endosulfan I	1	6.07	6.02	6.12	17.9	0
	2	5.25	5.20	5.30	17.9	
gamma-Chlordane	1	5.95	5.90	6.00	18.2	1.1
	2	5.13	5.08	5.18	18.4	
alpha-Chlordane	1	6.03	5.98	6.08	18.2	1.1
	2	5.19	5.14	5.24	18.0	
4,4'-DDE	1	6.20	6.15	6.25	18.0	0.6
	2	5.38	5.33	5.43	17.9	
Dieldrin	1	6.35	6.30	6.40	17.8	0.6
	2	5.51	5.46	5.56	17.9	
Endrin	1	6.57	6.52	6.62	16.9	1.8
	2	5.79	5.74	5.84	17.2	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

TP-76

Lab Name: Alliance

Contract: CAMP02

Lab Code: ACE

SDG NO.: Q2458

Lab Sample ID: Q2458-01

Date(s) Analyzed: 07/01/2025 07/01/2025

Instrument ID (1): ECD_D

Instrument ID (2): ECD_D

GC Column: (1): ZB-MR1

ID: 0.32 (mm)

GC Column:(2): ZB-MR2

ID: 0.32 (mm)

ANALYTE	COL	RT	RT WINDOW		CONCENTRATION	%RPD
			FROM	TO		
4,4'-DDT	1	7.03	6.98	7.08	0.51	80
	2	6.18	6.13	6.23	1.20	
Dieldrin	1	6.35	6.30	6.40	0.50	117.2
	2	5.51	5.46	5.56	1.90	
Endrin	1	6.58	6.53	6.63	0.66	8.1
	2	5.78	5.73	5.83	0.61	



QC SAMPLE DATA

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168672BL	SDG No.:	Q2458
Lab Sample ID:	PB168672BL	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	100 Decanted:
Sample Wt/Vol:	30.01 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089270.D	1	07/01/25 08:30	07/01/25 14:33	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	0.13	U	0.13	1.70	ug/kg
319-85-7	beta-BHC	0.18	U	0.18	1.70	ug/kg
319-86-8	delta-BHC	0.39	U	0.39	1.70	ug/kg
58-89-9	gamma-BHC (Lindane)	0.14	U	0.14	1.70	ug/kg
76-44-8	Heptachlor	0.12	U	0.12	1.70	ug/kg
309-00-2	Aldrin	0.12	U	0.12	1.70	ug/kg
1024-57-3	Heptachlor epoxide	0.19	U	0.19	1.70	ug/kg
959-98-8	Endosulfan I	0.14	U	0.14	1.70	ug/kg
60-57-1	Dieldrin	0.14	U	0.14	1.70	ug/kg
72-55-9	4,4-DDE	0.14	U	0.14	1.70	ug/kg
72-20-8	Endrin	0.14	U	0.14	1.70	ug/kg
33213-65-9	Endosulfan II	0.29	U	0.29	1.70	ug/kg
72-54-8	4,4-DDD	0.15	U	0.15	1.70	ug/kg
1031-07-8	Endosulfan Sulfate	0.13	U	0.13	1.70	ug/kg
50-29-3	4,4-DDT	0.14	U	0.14	1.70	ug/kg
72-43-5	Methoxychlor	0.37	U	0.37	1.70	ug/kg
53494-70-5	Endrin ketone	0.19	U	0.19	1.70	ug/kg
7421-93-4	Endrin aldehyde	0.37	U	0.37	1.70	ug/kg
5103-71-9	alpha-Chlordane	0.12	U	0.12	1.70	ug/kg
5103-74-2	gamma-Chlordane	0.15	U	0.15	1.70	ug/kg
8001-35-2	Toxaphene	5.40	U	5.40	33.0	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	17.3		20 - 144	87%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.0		19 - 148	100%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168672BL	SDG No.:	Q2458
Lab Sample ID:	PB168672BL	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	100 Decanted:
Sample Wt/Vol:	30.01 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089270.D	1	07/01/25 08:30	07/01/25 14:33	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089270.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 14:33
Operator : AR\AJ
Sample : PB168672BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PB168672BL

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:46:05 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	48153247	338.4E6	16.877	19.961
28)	SA Decachlor...	9.073	8.072	66591215	342.5E6	16.955	17.322

Target Compounds

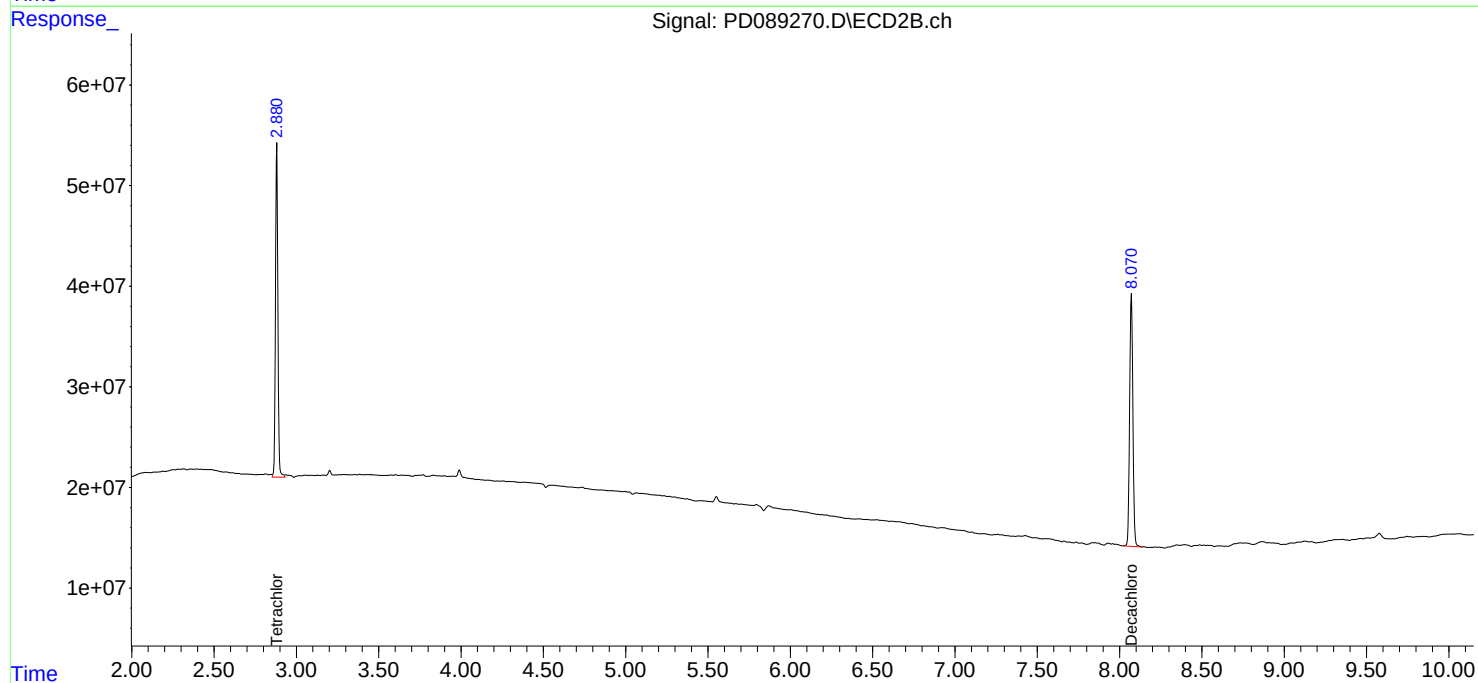
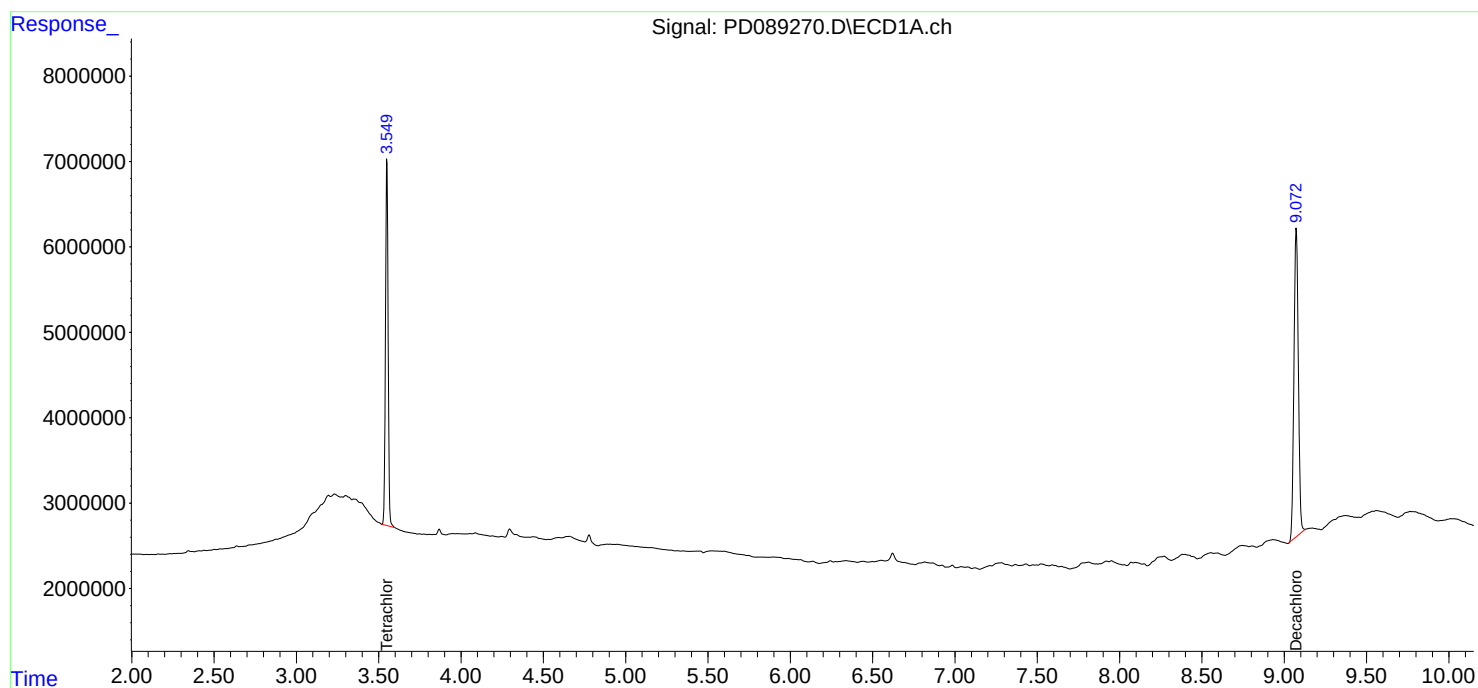
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

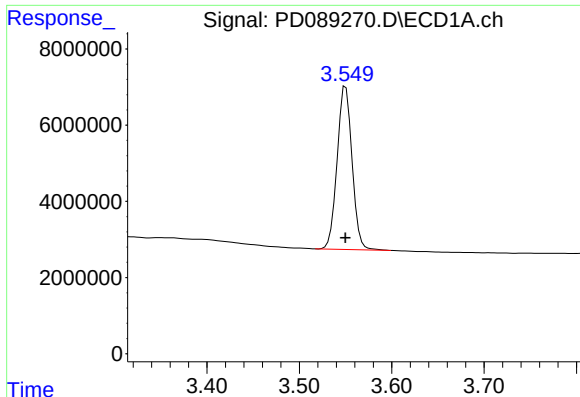
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089270.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 14:33
Operator : AR\AJ
Sample : PB168672BL
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PB168672BL

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:46:05 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

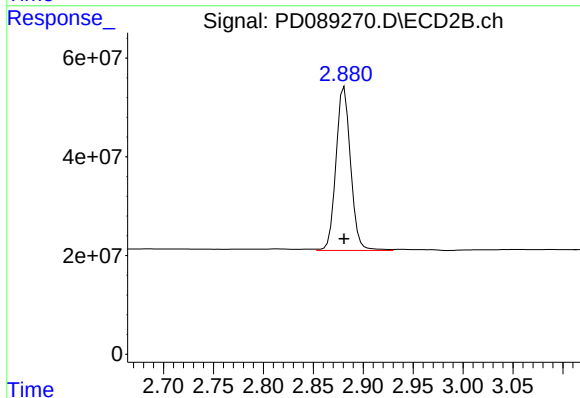




#1 Tetrachloro-m-xylene

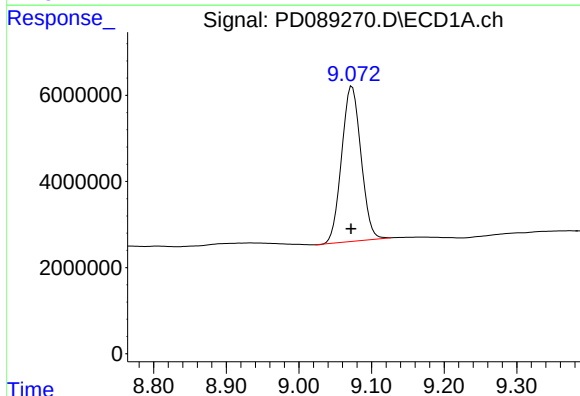
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 48153247
Conc: 16.88 ng/ml

Instrument :
ECD_D
ClientSampleId :
PB168672BL



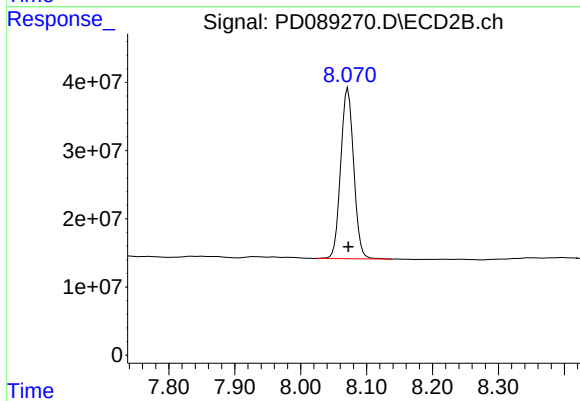
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 338405153
Conc: 19.96 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.001 min
Response: 66591215
Conc: 16.95 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 342462571
Conc: 17.32 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BL	SDG No.:	Q2458
Lab Sample ID:	PB168718BL	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089329.D	1	07/03/25 08:58	07/03/25 13:55	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	19.0		57 - 171	95%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.4		61 - 148	102%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BL	SDG No.:	Q2458
Lab Sample ID:	PB168718BL	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089329.D	1	07/03/25 08:58	07/03/25 13:55	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089329.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 13:55
 Operator : AR\AJ
 Sample : PB168718BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB168718BL

Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/07/2025
 Supervised By : mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:17:24 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds						
1) SA Tetrachlo...	3.557	2.879	52031894	345.5E6	18.237	20.379m
28) SA Decachlor...	9.084	8.075	73890119	376.5E6	18.813	19.044

Target Compounds

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089329.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 13:55
Operator : AR\AJ
Sample : PB168718BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

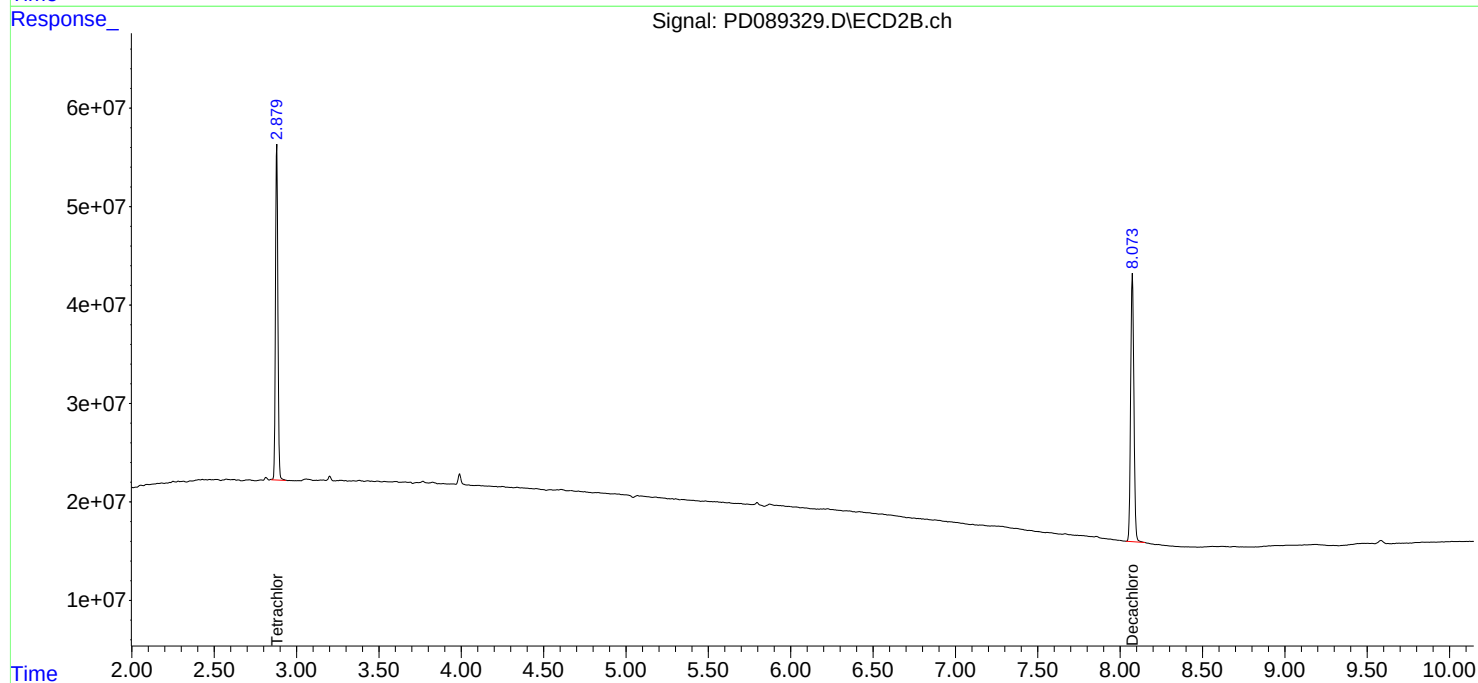
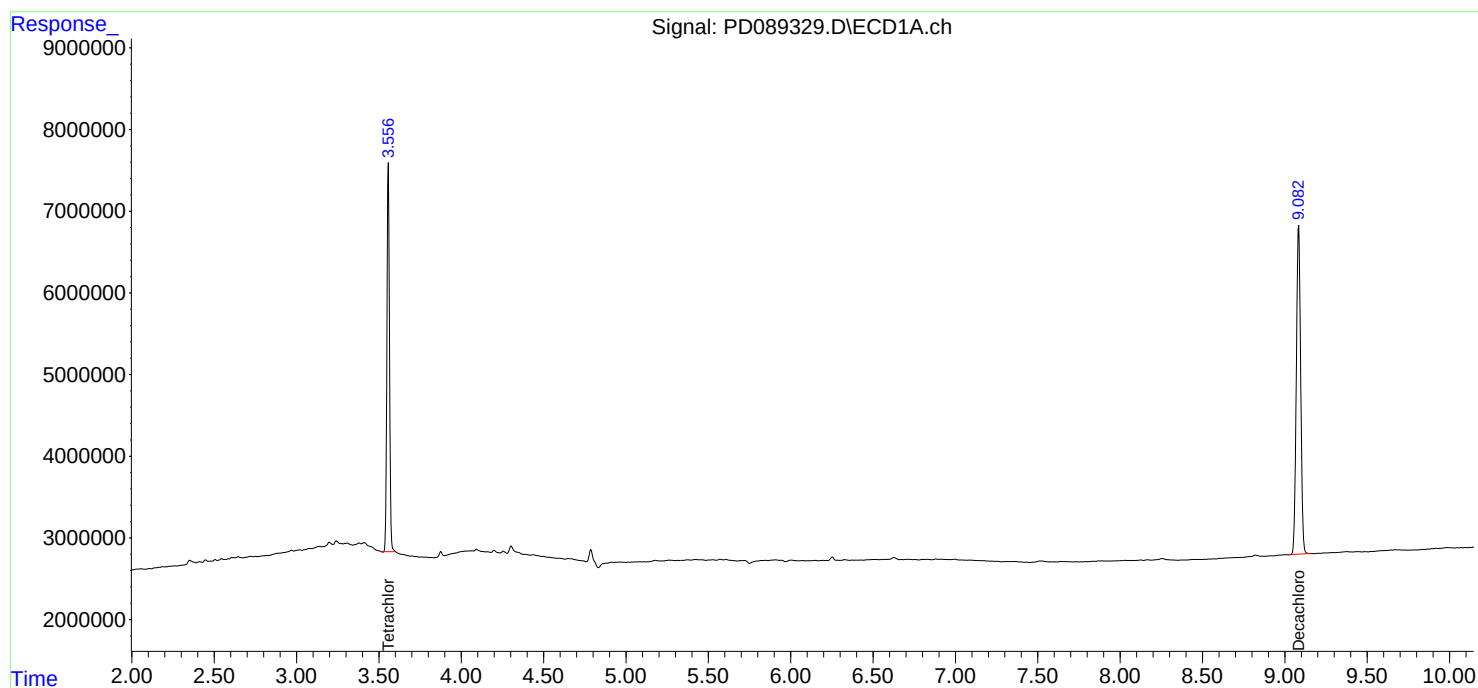
Instrument :
ECD_D
ClientSampleId :
PB168718BL

Manual Integrations
APPROVED

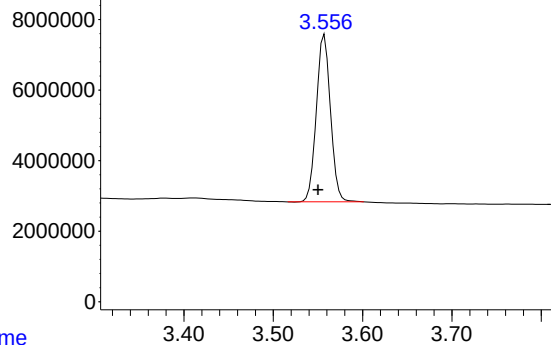
Reviewed By : Abdul Mirza 07/07/2025
Supervised By : mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:17:24 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Response_ Signal: PD089329.D\ECD1A.ch



#1 Tetrachloro-m-xylene

R.T.: 3.557 min
Delta R.T.: 0.007 min
Response: 52031894
Conc: 18.24 ng/ml

Instrument :

ECD_D

ClientSampleId :

PB168718BL

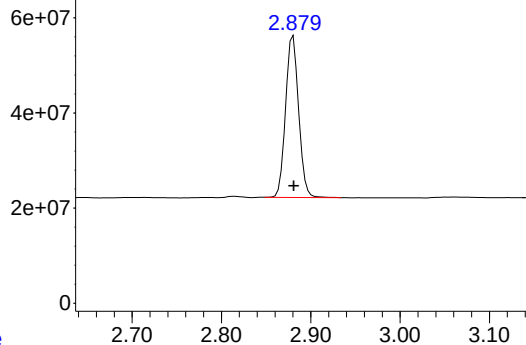
Manual Integrations

APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

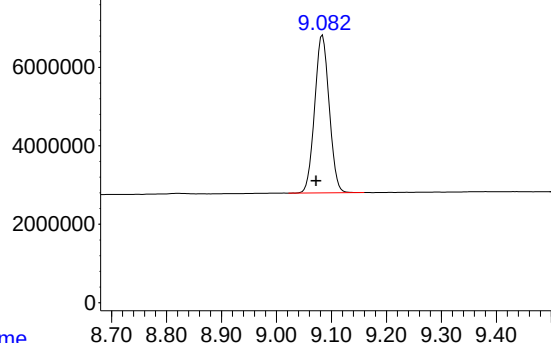
Time Response_ Signal: PD089329.D\ECD2B.ch



#1 Tetrachloro-m-xylene

R.T.: 2.879 min
Delta R.T.: -0.002 min
Response: 345485307
Conc: 20.38 ng/ml m

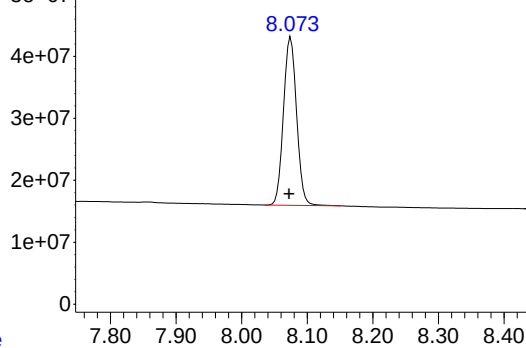
Time Response_ Signal: PD089329.D\ECD1A.ch



#28 Decachlorobiphenyl

R.T.: 9.084 min
Delta R.T.: 0.012 min
Response: 73890119
Conc: 18.81 ng/ml

Time Response_ Signal: PD089329.D\ECD2B.ch



#28 Decachlorobiphenyl

R.T.: 8.075 min
Delta R.T.: 0.003 min
Response: 376506282
Conc: 19.04 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	06/17/25
Project:	South River WM Replacement	Date Received:	06/17/25
Client Sample ID:	PIBLK-PD088990.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD088990.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD088990.D	1		06/17/25	PD061825

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	18.6		57 - 171	93%	SPK: 20
877-09-8	Tetrachloro-m-xylene	17.7		61 - 148	88%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/17/25
Project:	South River WM Replacement	Date Received:	06/17/25
Client Sample ID:	PIBLK-PD088990.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD088990.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD088990.D	1		06/17/25	PD061825

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088990.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 15:11
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:45:07 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.549	2.881	45828945	299.5E6	16.063	17.668
28)	SA Decachlor...	9.072	8.072	71428242	368.1E6	18.186	18.617

Target Compounds

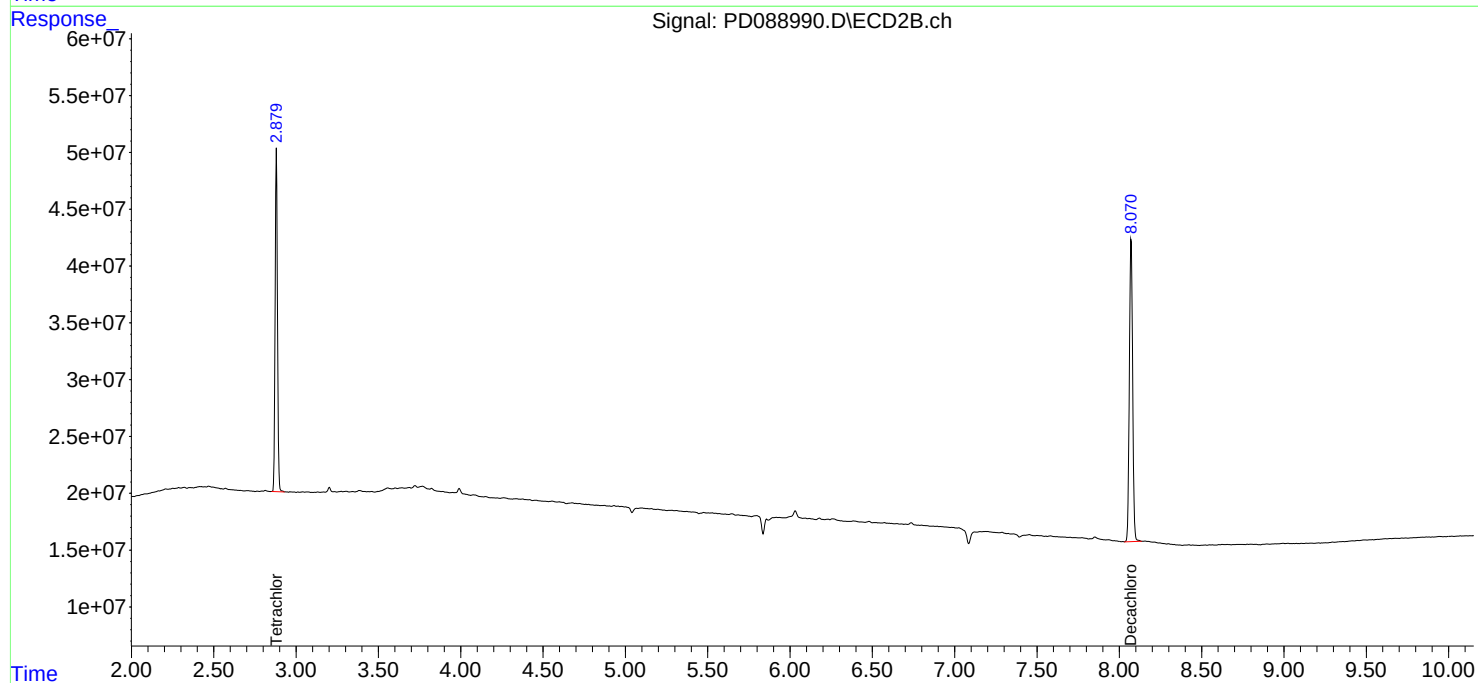
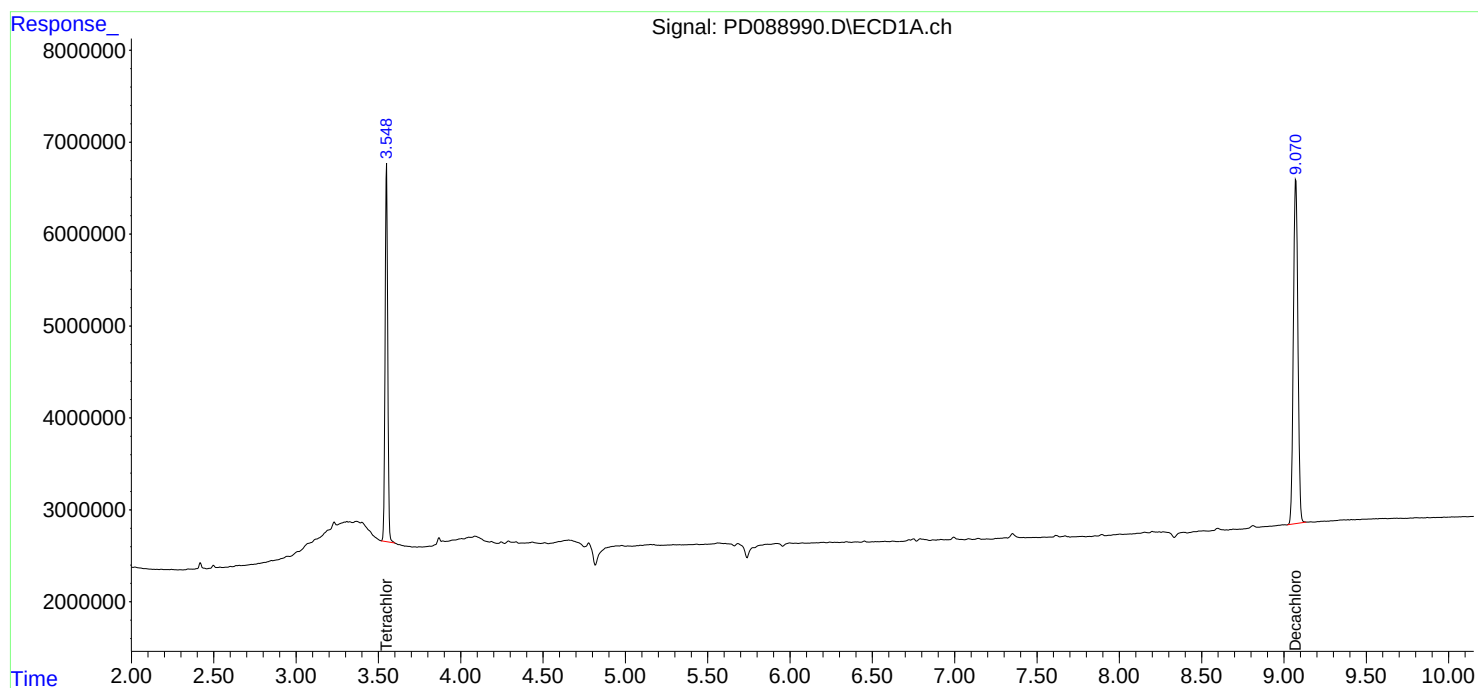
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

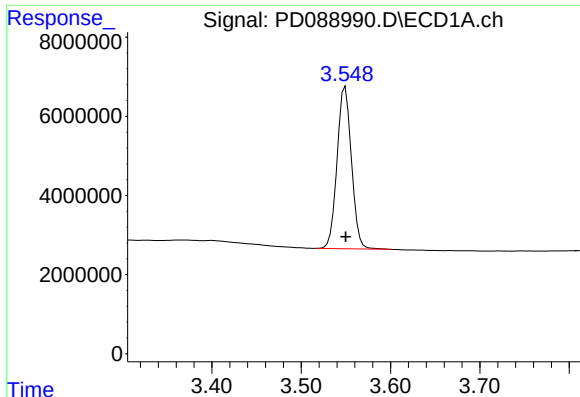
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD061825\
Data File : PD088990.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 17 Jun 2025 15:11
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jun 18 05:45:07 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

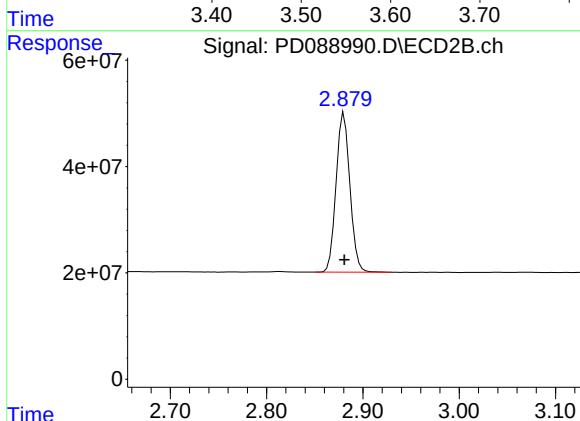




#1 Tetrachloro-m-xylene

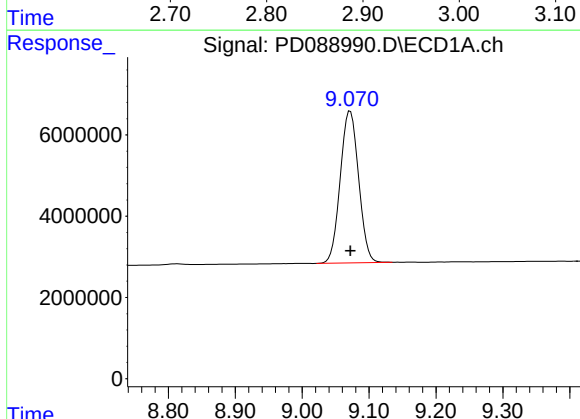
R.T.: 3.549 min
Delta R.T.: 0.000 min
Response: 45828945
Conc: 16.06 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



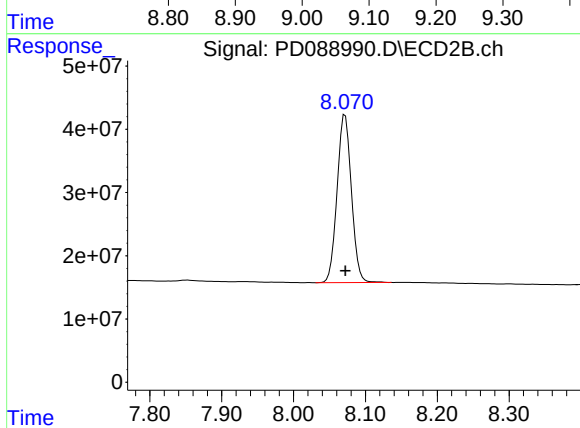
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 299529104
Conc: 17.67 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.072 min
Delta R.T.: 0.000 min
Response: 71428242
Conc: 18.19 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 368061815
Conc: 18.62 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089264.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089264.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089264.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	23.3		57 - 171	117%	SPK: 20
877-09-8	Tetrachloro-m-xylene	22.7		61 - 148	114%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089264.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089264.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089264.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089264.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 11:59
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:44:38 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.881	58707613	385.4E6	20.577	22.732
28)	SA Decachlor...	9.073	8.072	89993686	460.9E6	22.913	23.314

Target Compounds

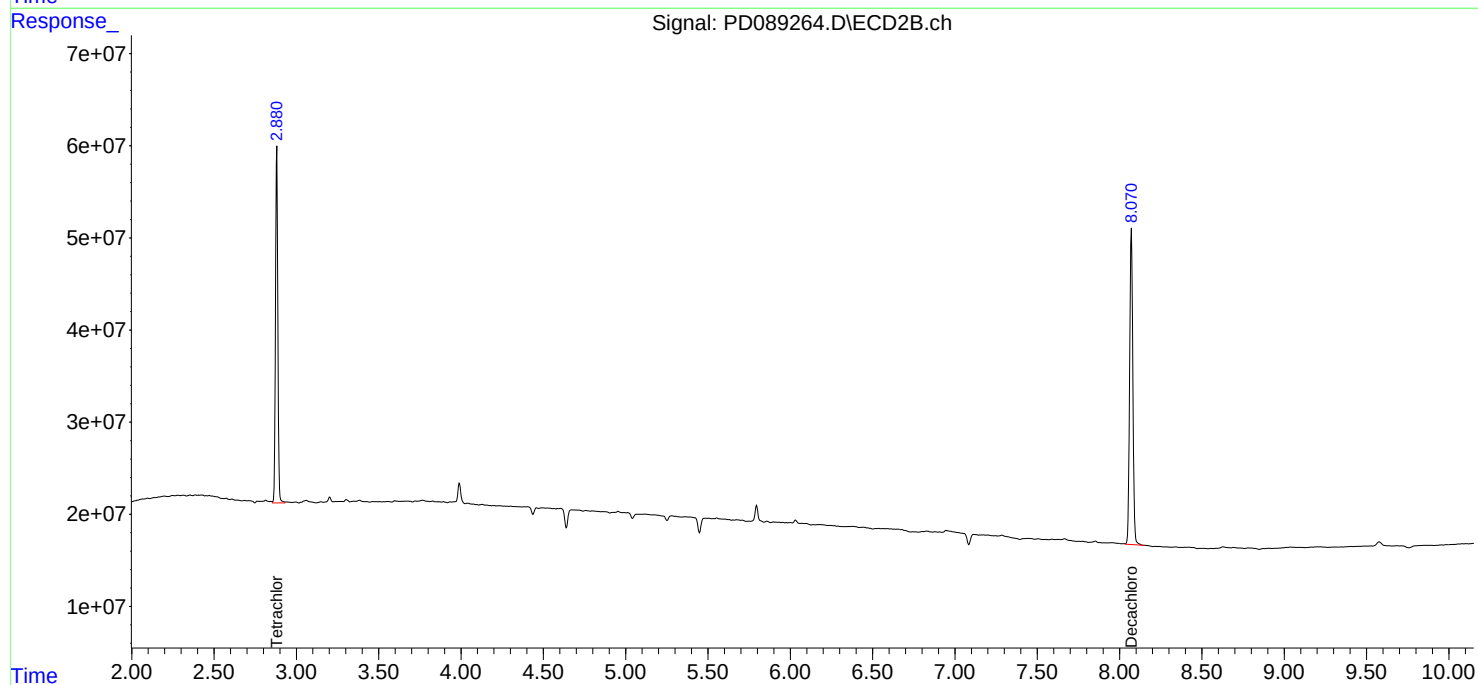
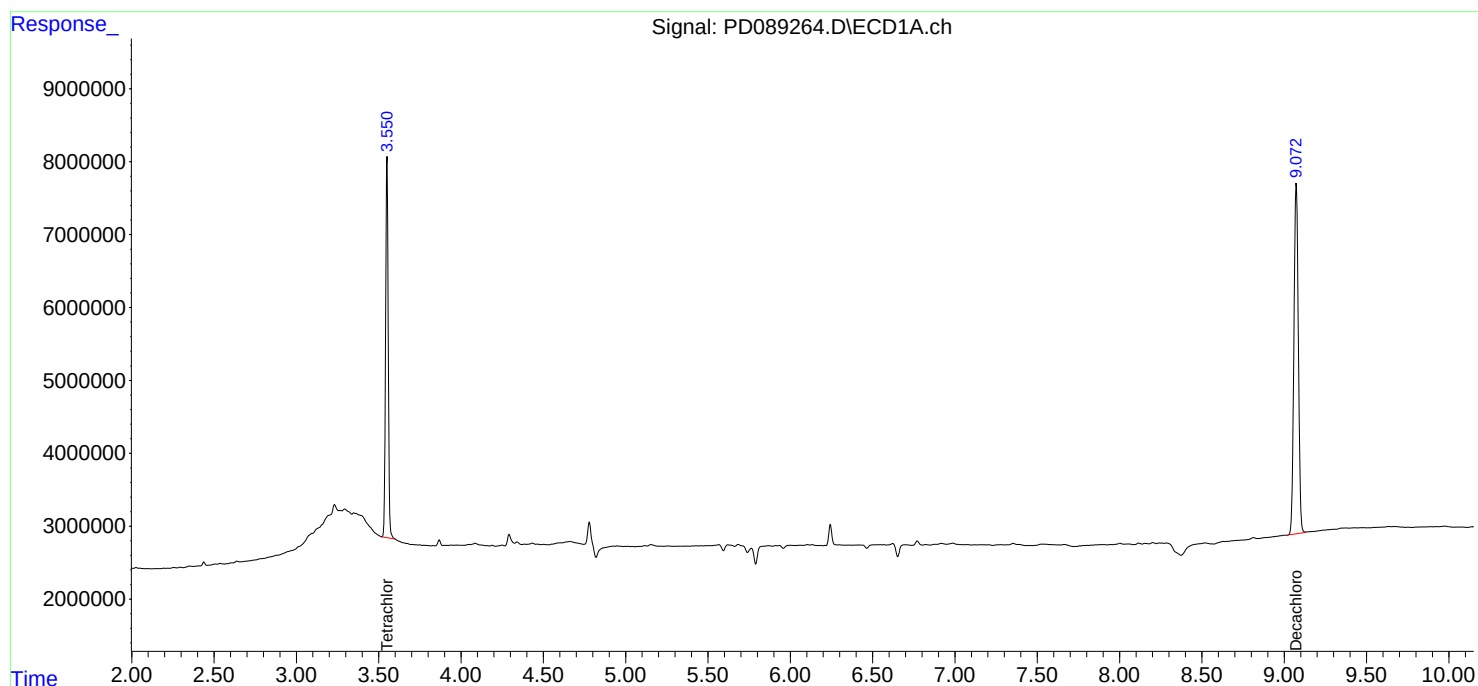
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

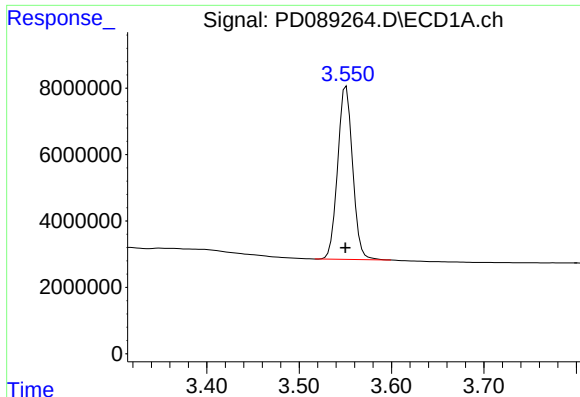
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089264.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 11:59
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:44:38 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

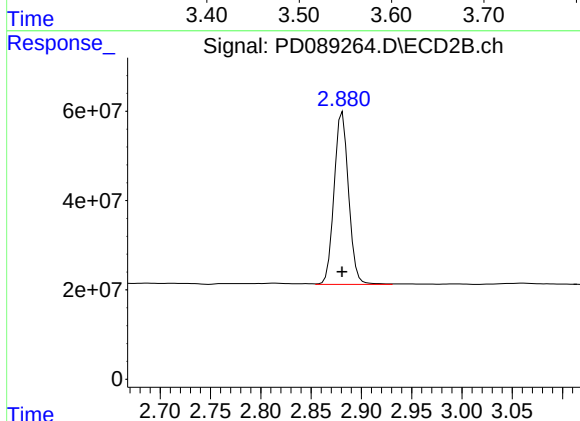




#1 Tetrachloro-m-xylene

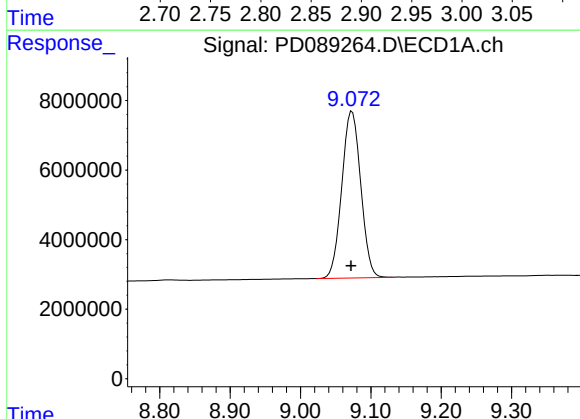
R.T.: 3.551 min
Delta R.T.: 0.001 min
Response: 58707613
Conc: 20.58 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



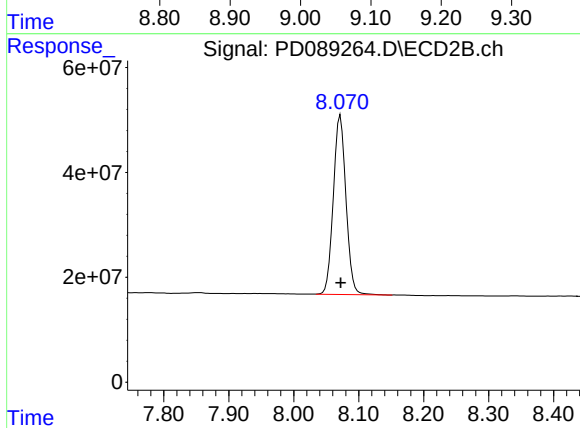
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 385371808
Conc: 22.73 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.073 min
Delta R.T.: 0.001 min
Response: 89993686
Conc: 22.91 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 460913349
Conc: 23.31 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089276.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089276.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089276.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	23.0		57 - 171	115%	SPK: 20
877-09-8	Tetrachloro-m-xylene	23.2		61 - 148	116%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089276.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089276.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089276.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089276.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 16:37
 Operator : AR\AJ
 Sample : I.BLK
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 I.BLK

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:47:46 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	59714817	392.7E6	20.930	23.165
28)	SA Decachlor...	9.074	8.072	90269419	443.3E6	22.983	22.421

Target Compounds

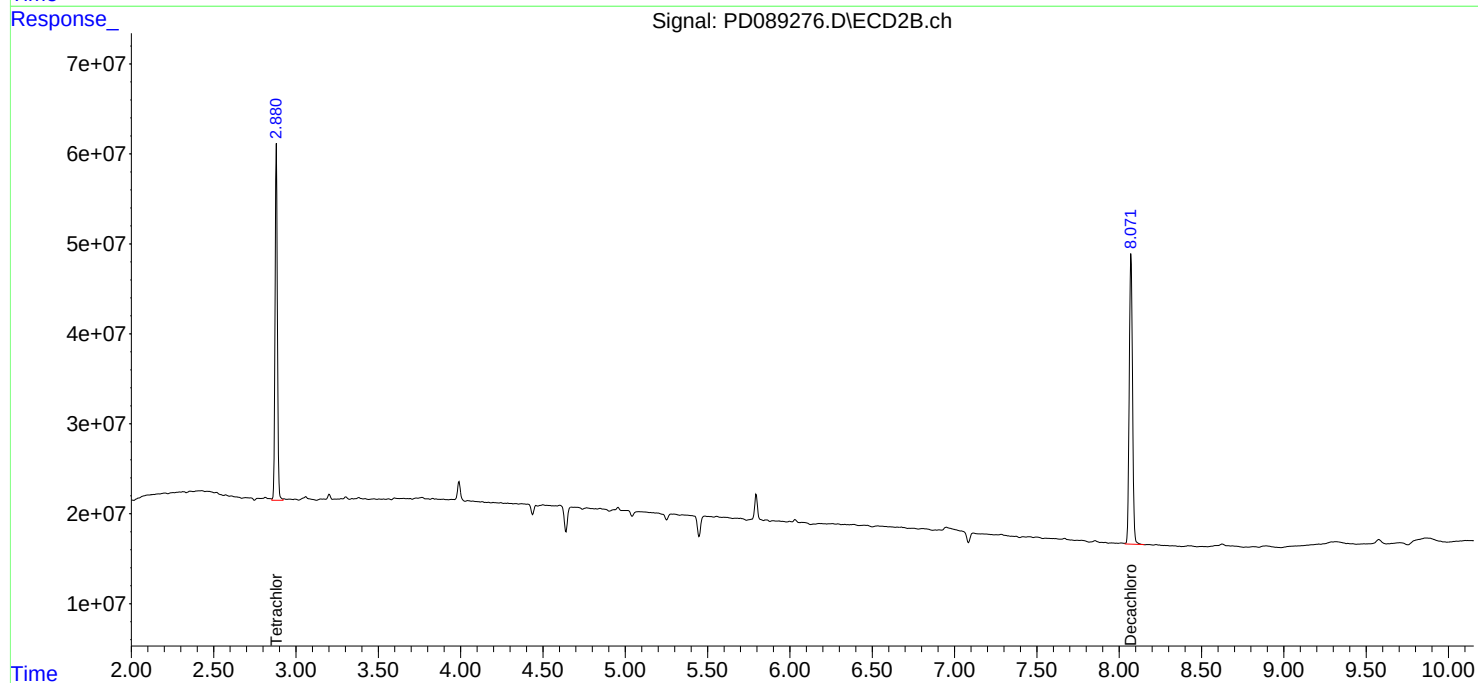
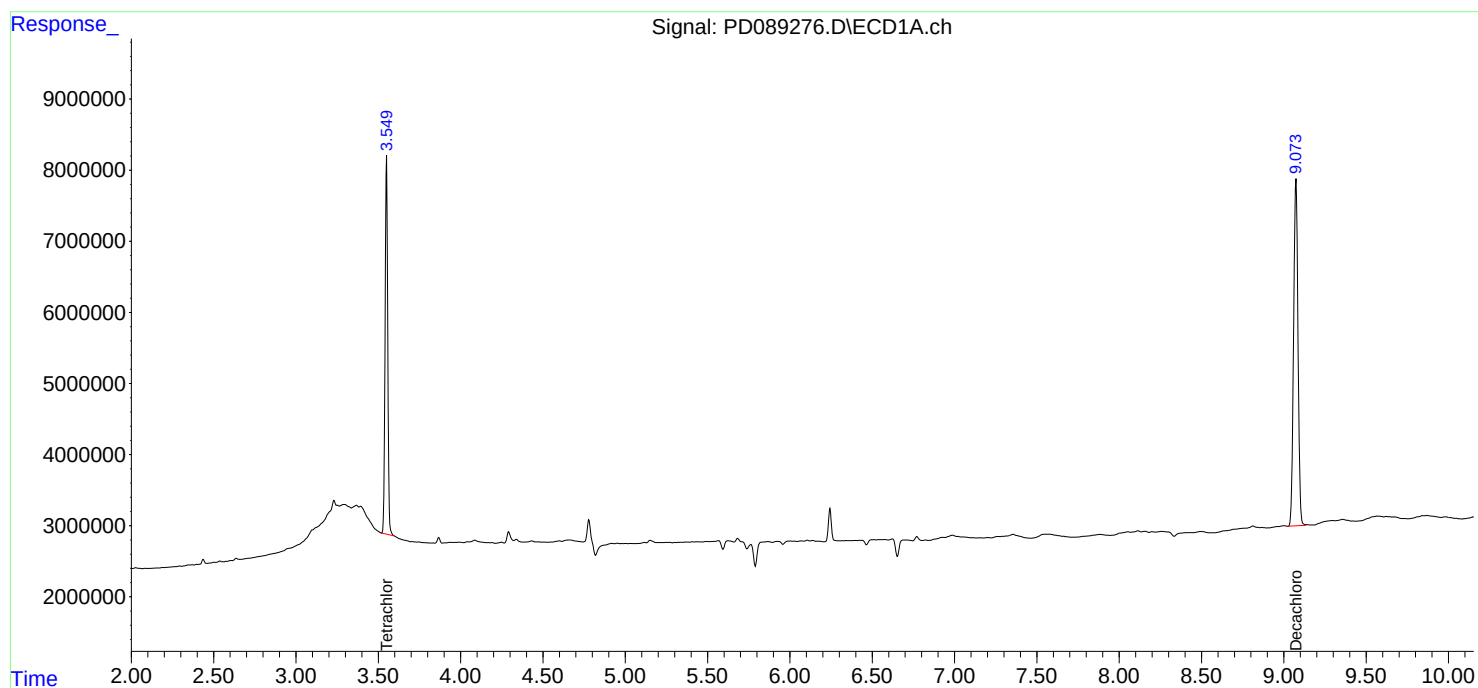
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

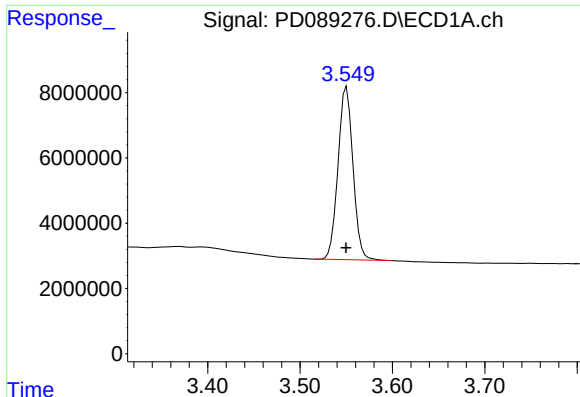
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089276.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 16:37
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:47:46 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

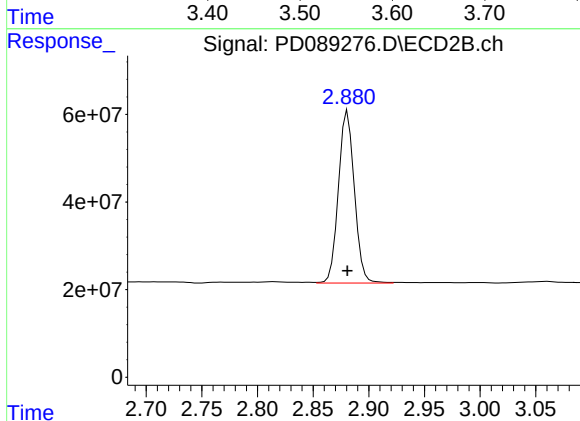




#1 Tetrachloro-m-xylene

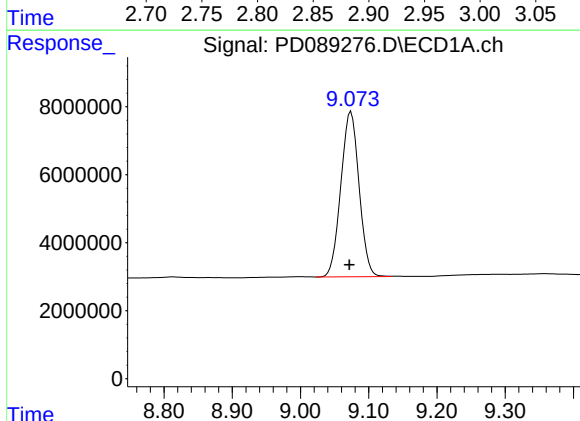
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 59714817
Conc: 20.93 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



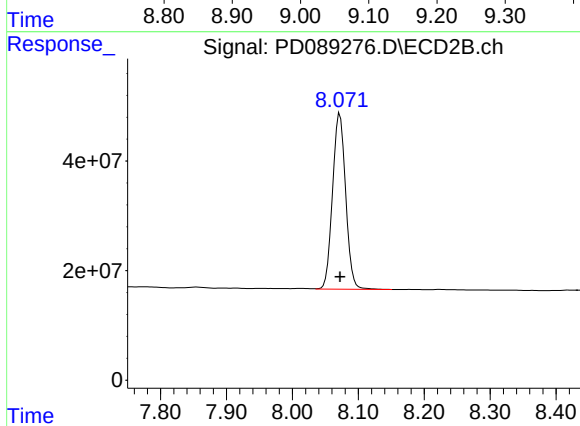
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 392715775
Conc: 23.16 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.074 min
Delta R.T.: 0.002 min
Response: 90269419
Conc: 22.98 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 443269881
Conc: 22.42 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089287.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089287.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089287.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	22.1		57 - 171	110%	SPK: 20
877-09-8	Tetrachloro-m-xylene	23.4		61 - 148	117%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089287.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089287.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089287.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089287.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 19:07
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:50:23 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	59834557	396.9E6	20.972	23.413
28)	SA Decachlor...	9.072	8.071	86680293	396.6E6	22.069	20.060

Target Compounds

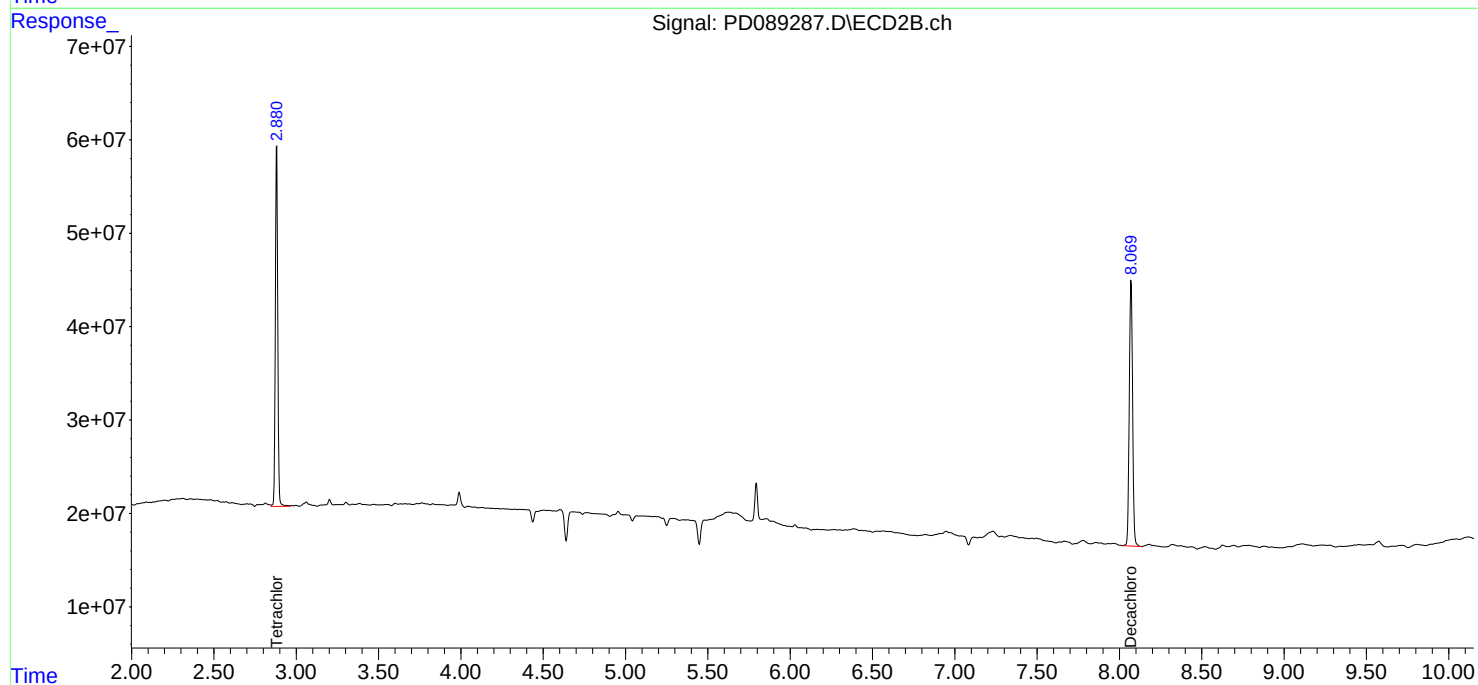
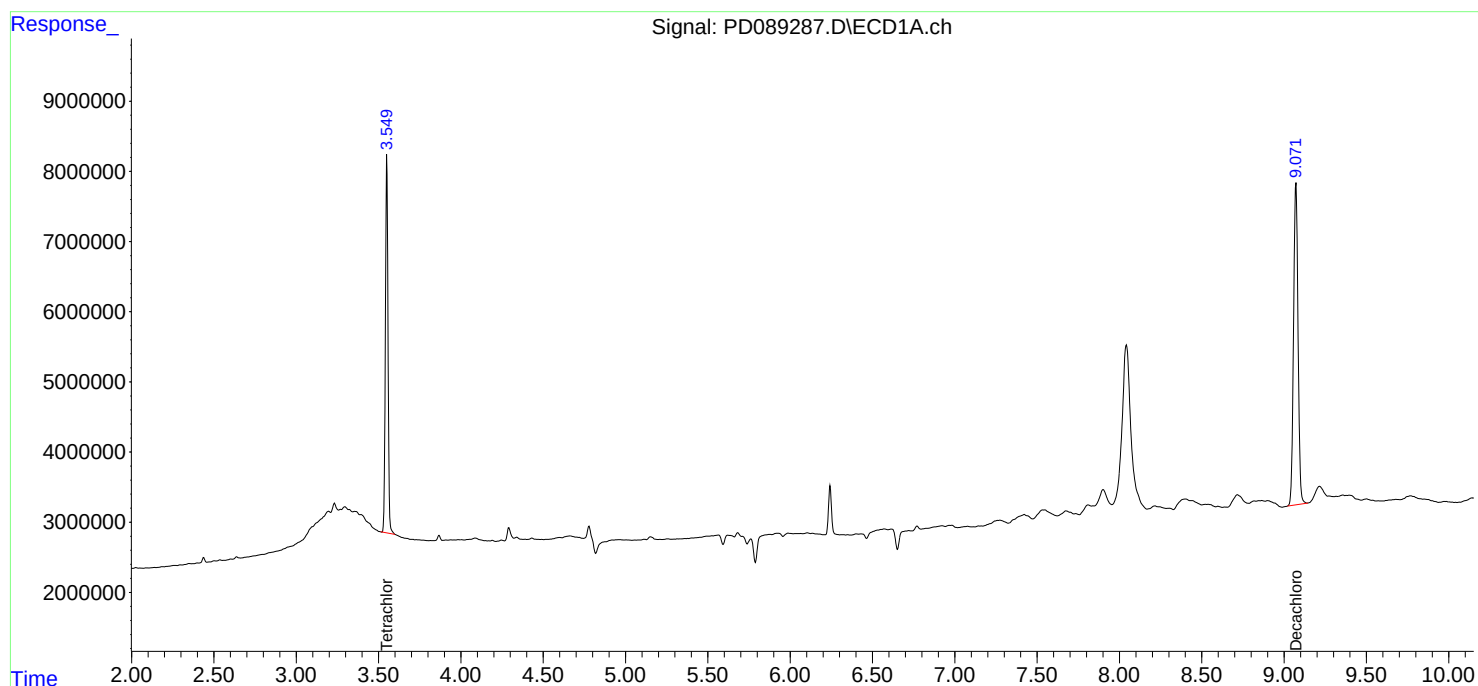
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

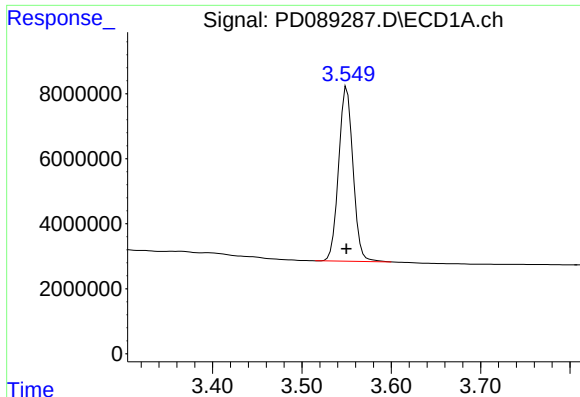
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089287.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 19:07
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:50:23 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

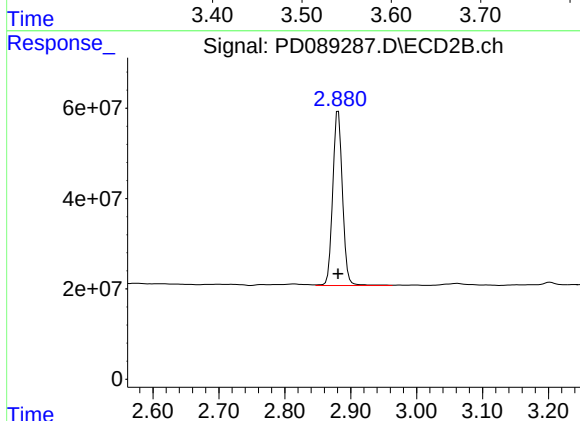




#1 Tetrachloro-m-xylene

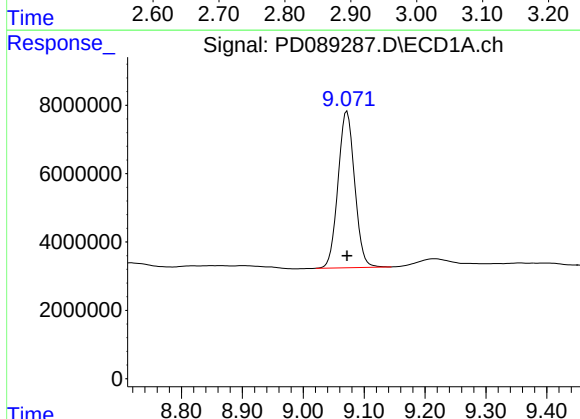
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 59834557
Conc: 20.97 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



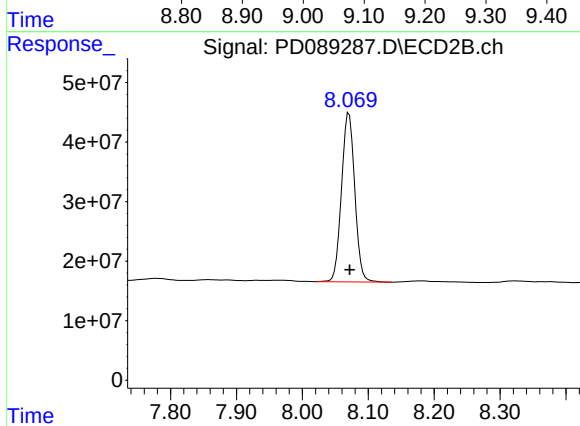
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 396926169
Conc: 23.41 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.072 min
Delta R.T.: 0.000 min
Response: 86680293
Conc: 22.07 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.071 min
Delta R.T.: -0.001 min
Response: 396587260
Conc: 20.06 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089298.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089298.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089298.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	21.1		57 - 171	106%	SPK: 20
877-09-8	Tetrachloro-m-xylene	24.2		61 - 148	121%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/01/25
Project:	South River WM Replacement	Date Received:	07/01/25
Client Sample ID:	PIBLK-PD089298.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089298.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089298.D	1		07/01/25	pd070225

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089298.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 22:05
 Operator : AR\AJ
 Sample : I.BLK
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 I.BLK

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:52:59 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	61643073	410.3E6	21.605	24.204
28)	SA Decachlor...	9.072	8.072	83023185	359.7E6	21.138	18.195

Target Compounds

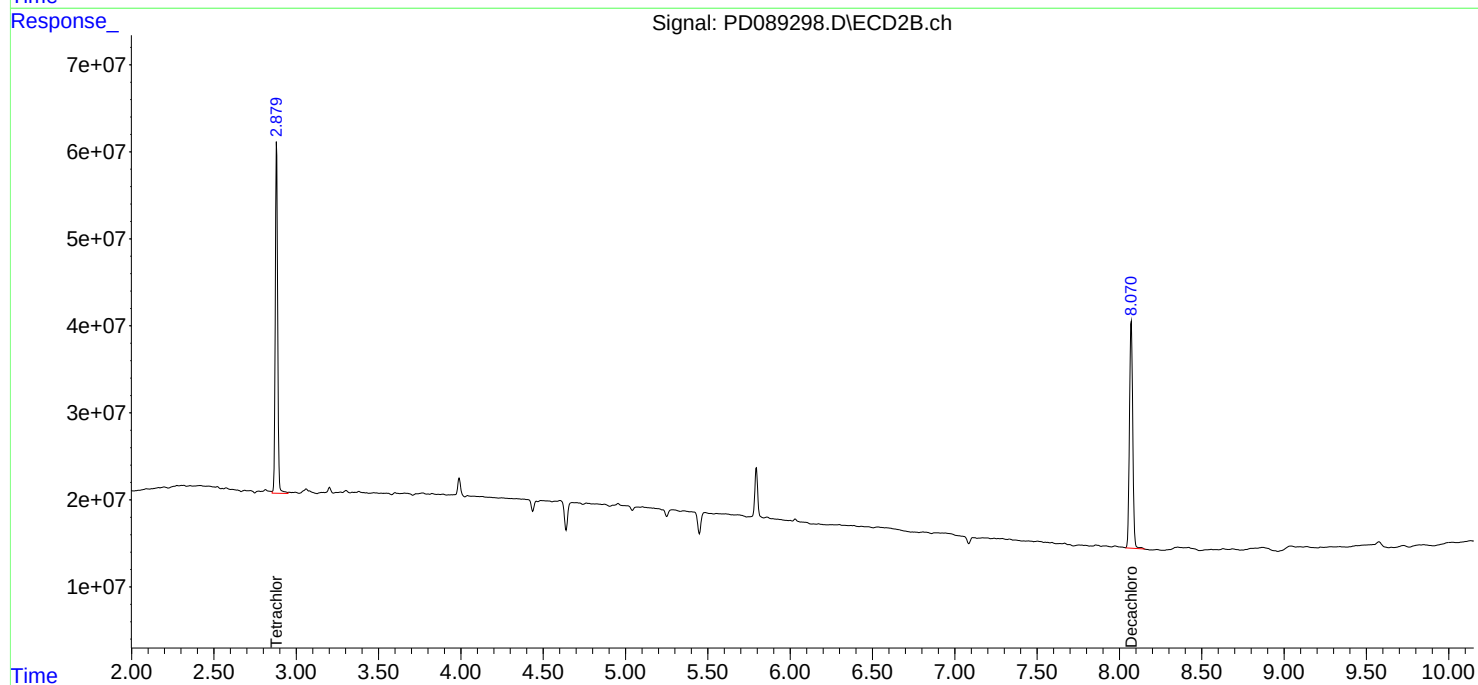
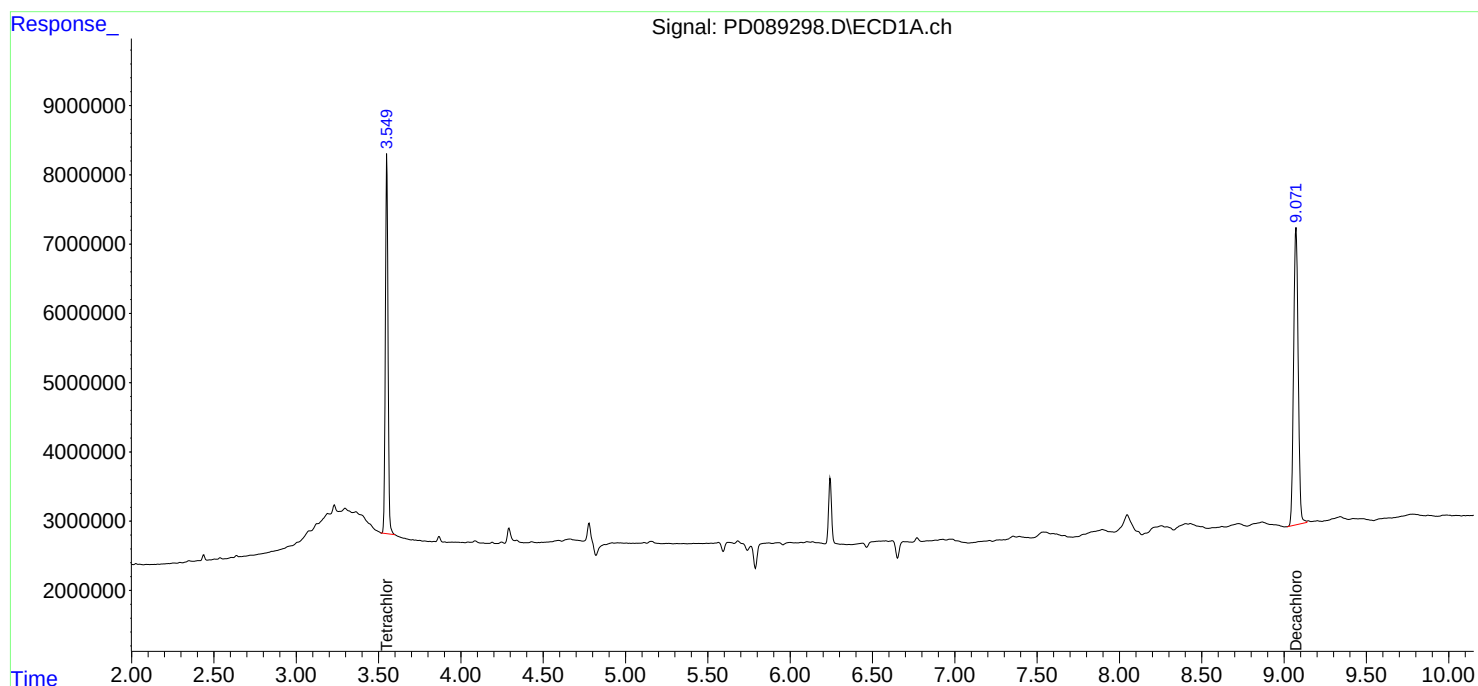
 (f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

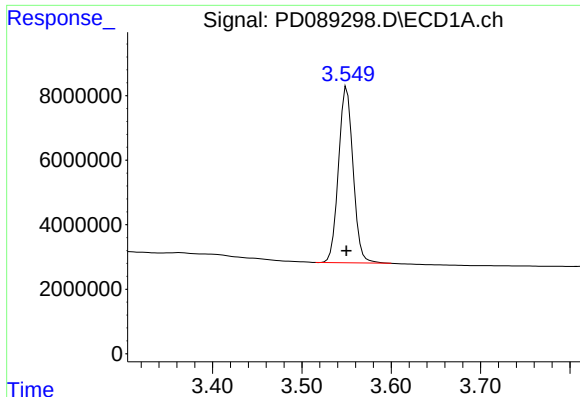
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089298.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 22:05
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:52:59 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

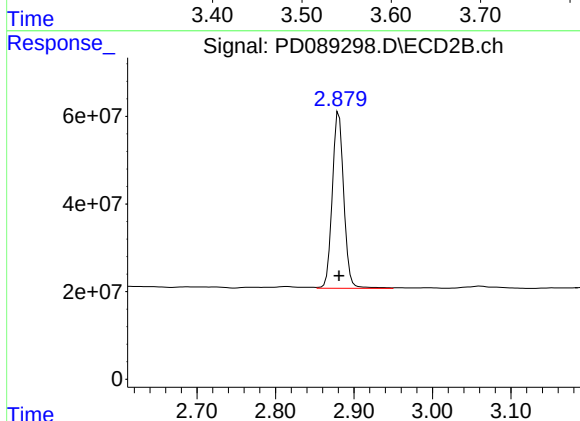




#1 Tetrachloro-m-xylene

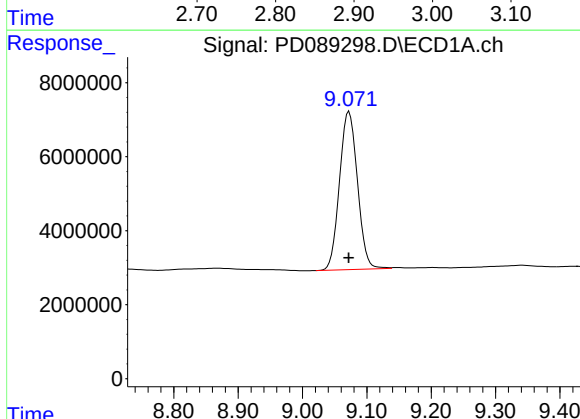
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 61643073
Conc: 21.61 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



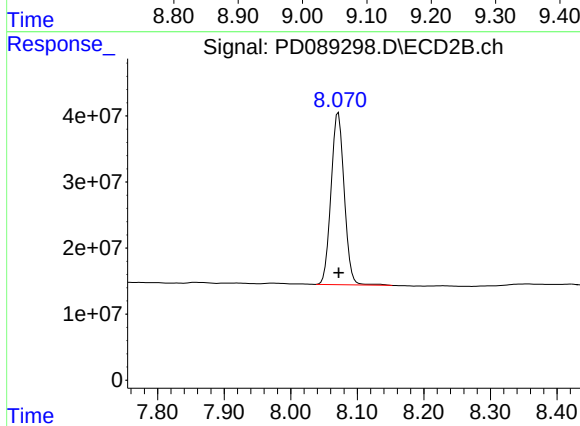
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 410342837
Conc: 24.20 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.072 min
Delta R.T.: 0.000 min
Response: 83023185
Conc: 21.14 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 359720322
Conc: 18.20 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/03/25
Project:	South River WM Replacement	Date Received:	07/03/25
Client Sample ID:	PIBLK-PD089326.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089326.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089326.D	1		07/03/25	pd070425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	19.1		57 - 171	96%	SPK: 20
877-09-8	Tetrachloro-m-xylene	21.9		61 - 148	109%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/03/25
Project:	South River WM Replacement	Date Received:	07/03/25
Client Sample ID:	PIBLK-PD089326.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089326.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089326.D	1		07/03/25	pd070425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089326.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 09:19
 Operator : AR\AJ
 Sample : I.BLK
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 I.BLK

Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/07/2025
 Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:16:37 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.879	55547175	371.1E6	19.469	21.890m
28)	SA Decachlor...	9.076	8.071	75117817	355.5E6	19.125	17.980

Target Compounds

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089326.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 09:19
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

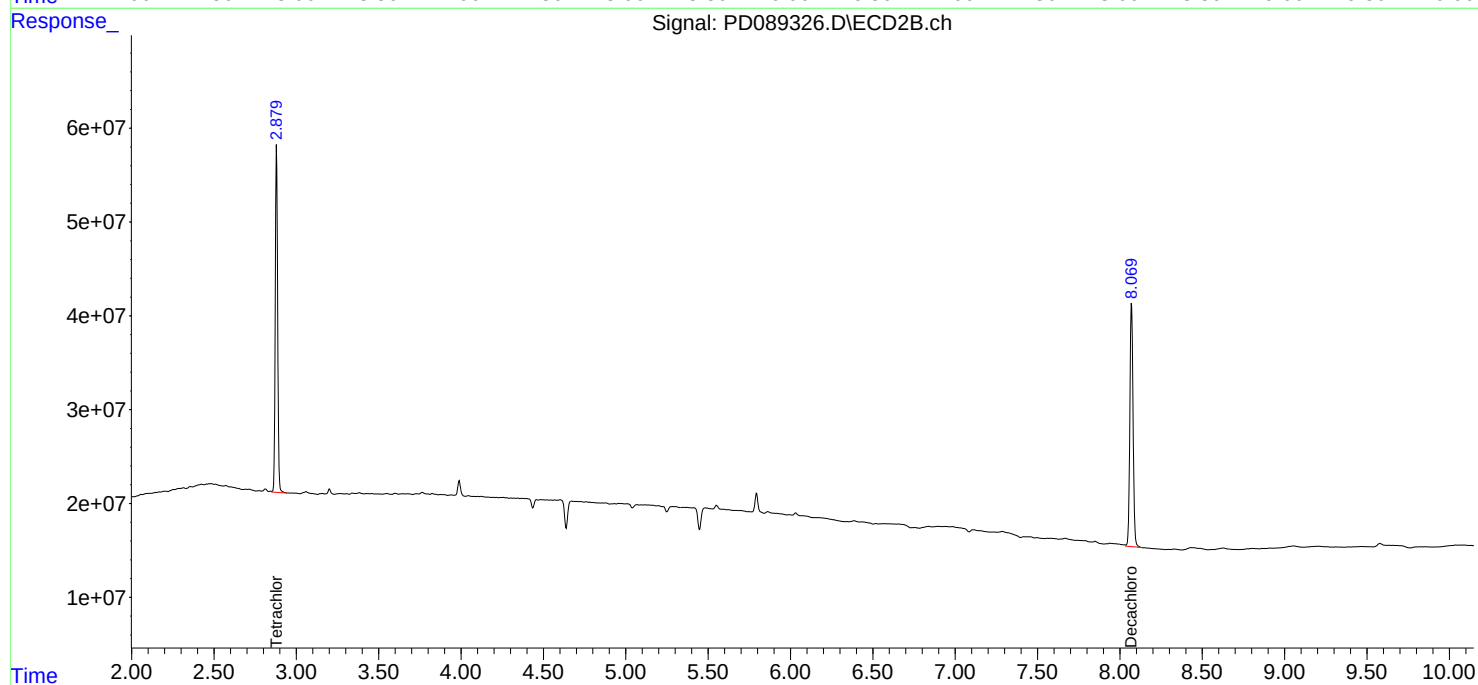
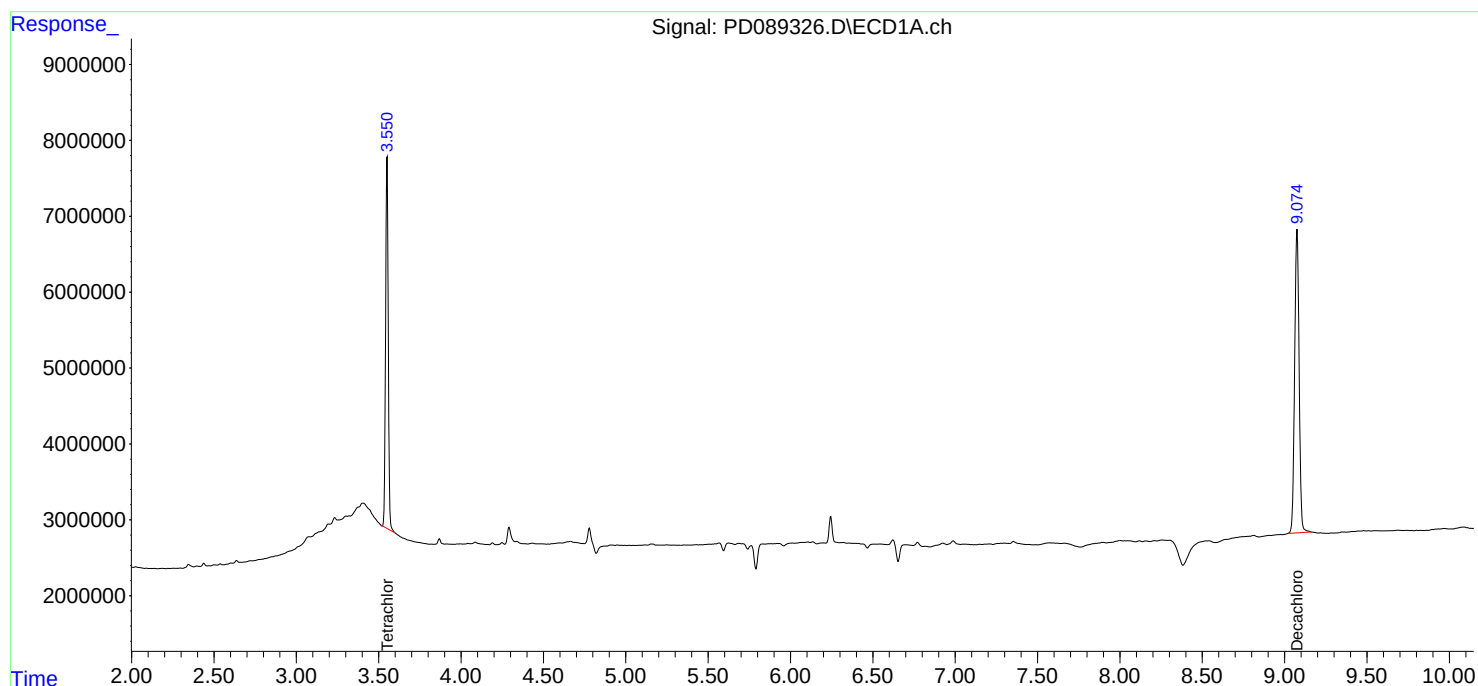
Instrument :
ECD_D
ClientSampleId :
I.BLK

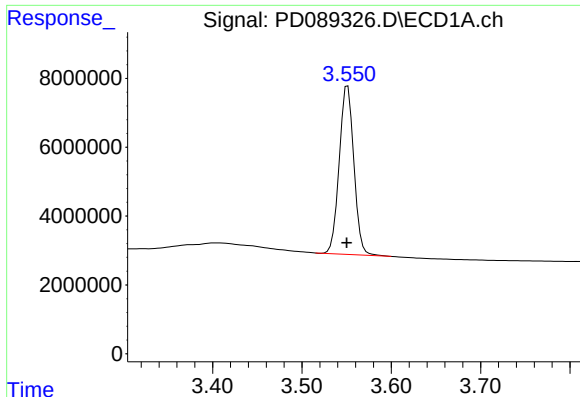
Manual Integrations
APPROVED

Reviewed By : Abdul Mirza 07/07/2025
Supervised By : mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:16:37 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





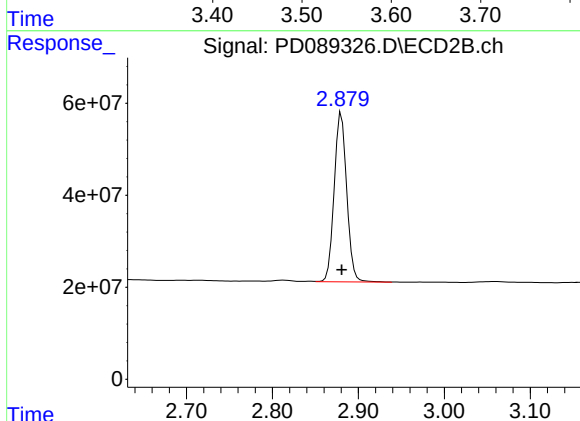
#1 Tetrachloro-m-xylene

R.T.: 3.551 min
Delta R.T.: 0.001 min
Response: 55547175
Conc: 19.47 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK

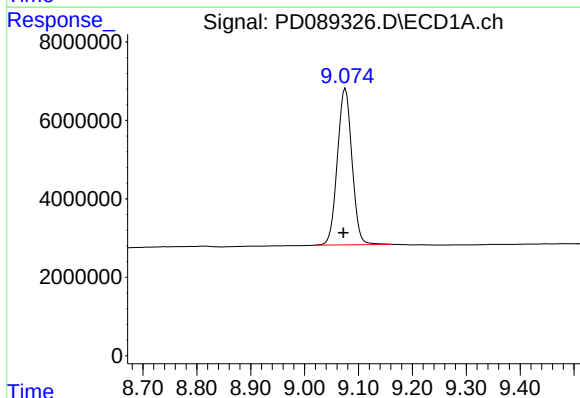
Manual Integrations
APPROVED

Reviewed By :Abdul Mirza 07/07/2025
Supervised By :mohammad ahmed 07/09/2025



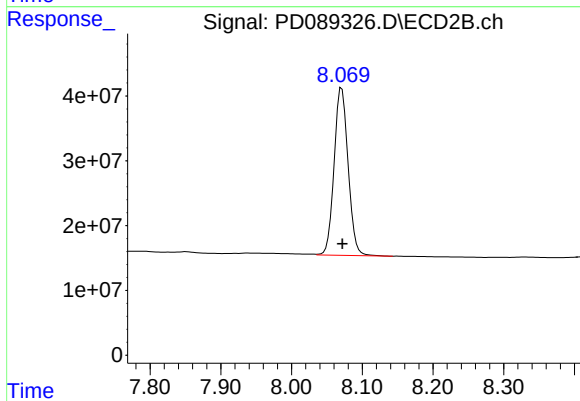
#1 Tetrachloro-m-xylene

R.T.: 2.879 min
Delta R.T.: -0.002 min
Response: 371105420
Conc: 21.89 ng/ml m



#28 Decachlorobiphenyl

R.T.: 9.076 min
Delta R.T.: 0.004 min
Response: 75117817
Conc: 19.13 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.071 min
Delta R.T.: -0.001 min
Response: 355472700
Conc: 17.98 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	07/03/25
Project:	South River WM Replacement	Date Received:	07/03/25
Client Sample ID:	PIBLK-PD089337.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089337.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089337.D	1		07/03/25	pd070425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.0039	U	0.0039	0.050	ug/L
319-85-7	beta-BHC	0.0049	U	0.0049	0.050	ug/L
319-86-8	delta-BHC	0.011	U	0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.0037	U	0.0037	0.050	ug/L
76-44-8	Heptachlor	0.0027	U	0.0027	0.050	ug/L
309-00-2	Aldrin	0.0036	U	0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.0096	U	0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.0031	U	0.0031	0.050	ug/L
60-57-1	Dieldrin	0.0036	U	0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.0037	U	0.0037	0.050	ug/L
72-20-8	Endrin	0.0032	U	0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.0079	U	0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.0071	U	0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.0037	U	0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.0035	U	0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.011	U	0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.0093	U	0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.011	U	0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.0035	U	0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.0039	U	0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	19.1		57 - 171	95%	SPK: 20
877-09-8	Tetrachloro-m-xylene	22.4		61 - 148	112%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	07/03/25
Project:	South River WM Replacement	Date Received:	07/03/25
Client Sample ID:	PIBLK-PD089337.D	SDG No.:	Q2458
Lab Sample ID:	I.BLK-PD089337.D	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0
Sample Wt/Vol:	1000	Units:	mL
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	PH :	
Prep Method :	3510C	Injection Volume :	

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089337.D	1		07/03/25	pd070425

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089337.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 15:48
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:19:28 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.881	56361551	379.8E6	19.754	22.405
28)	SA Decachlor...	9.075	8.072	74529740	377.3E6	18.976	19.086

Target Compounds

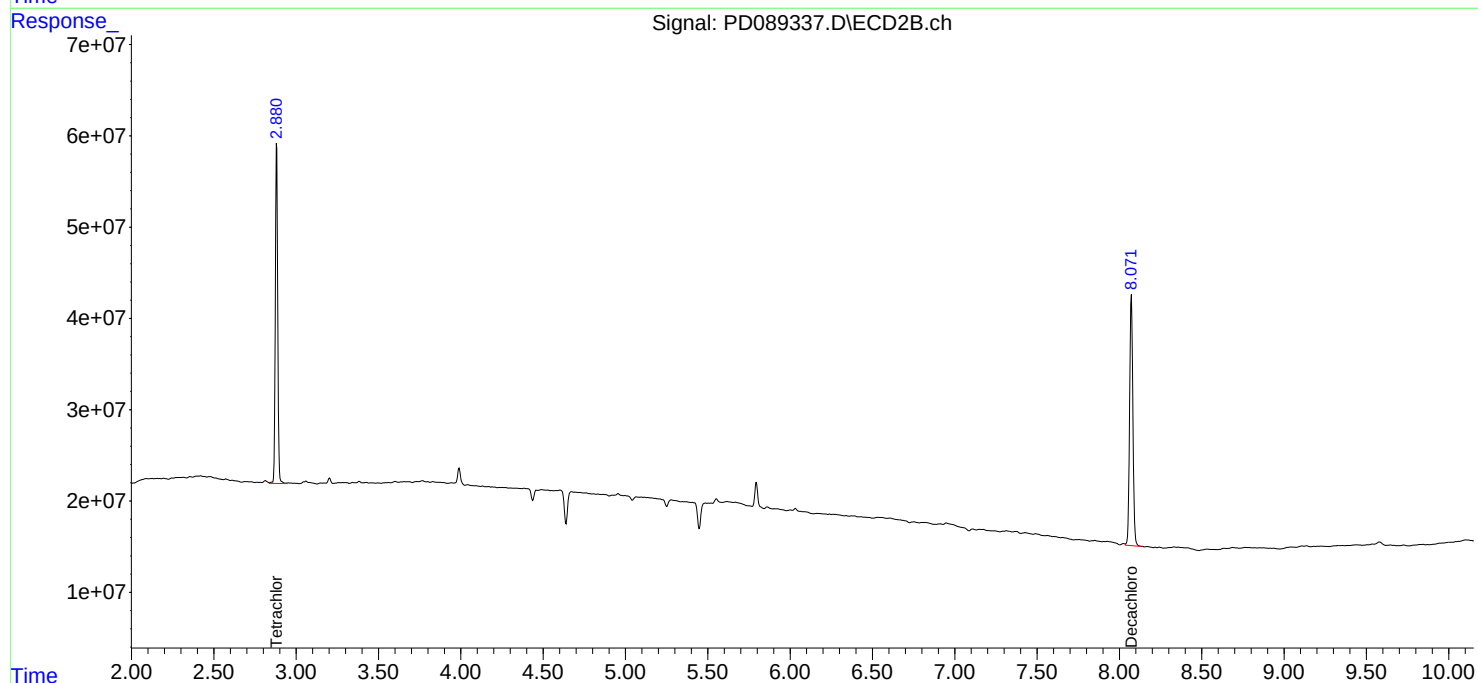
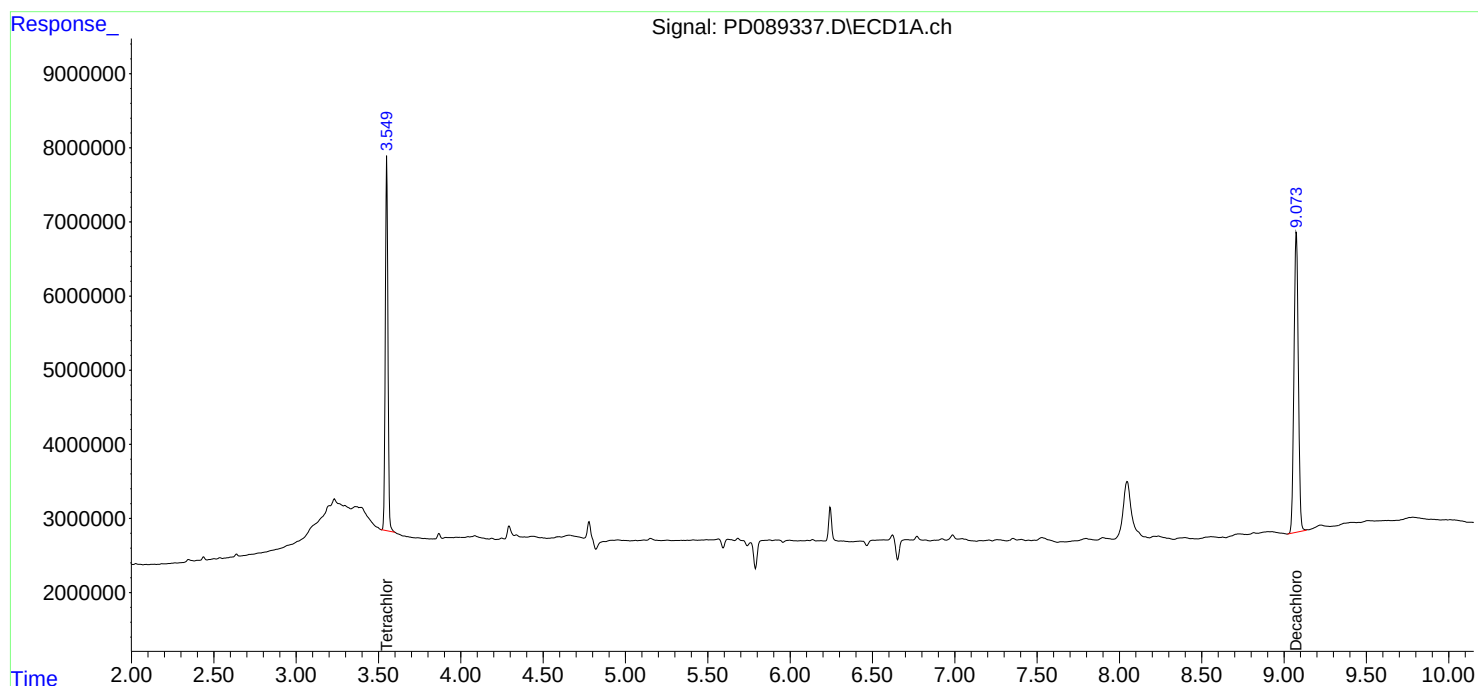
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

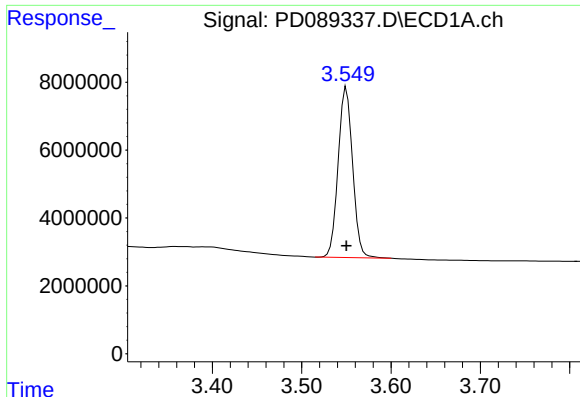
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089337.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 15:48
Operator : AR\AJ
Sample : I.BLK
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
I.BLK

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:19:28 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

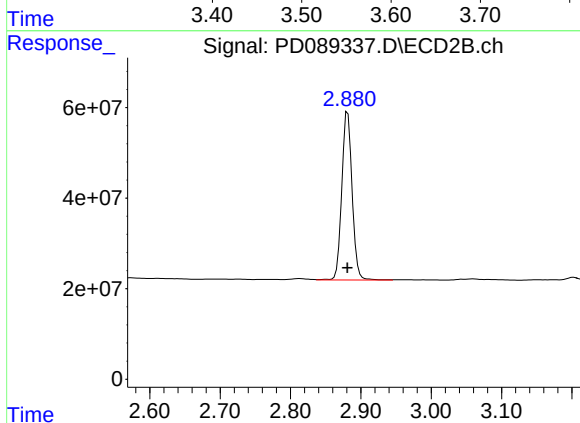




#1 Tetrachloro-m-xylene

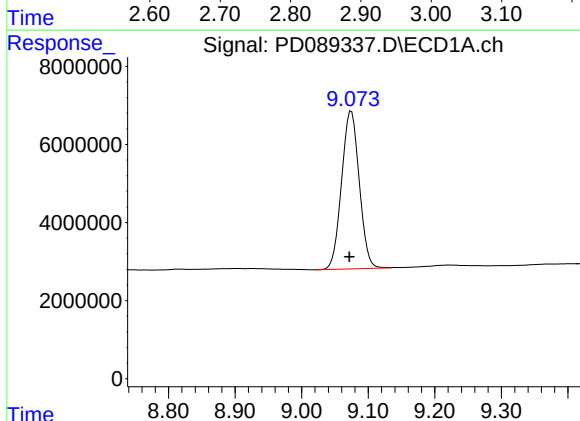
R.T.: 3.550 min
Delta R.T.: 0.000 min
Response: 56361551
Conc: 19.75 ng/ml

Instrument :
ECD_D
ClientSampleId :
I.BLK



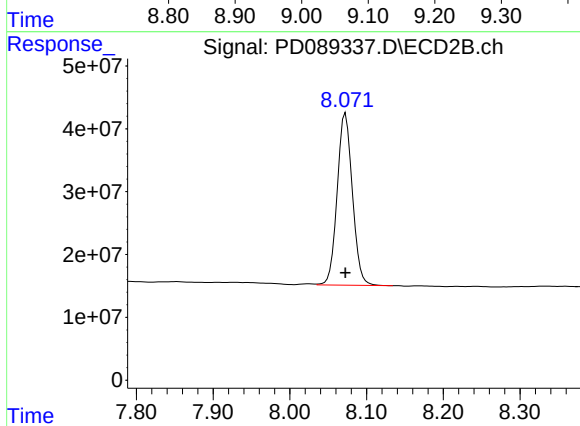
#1 Tetrachloro-m-xylene

R.T.: 2.881 min
Delta R.T.: 0.000 min
Response: 379829840
Conc: 22.40 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.075 min
Delta R.T.: 0.003 min
Response: 74529740
Conc: 18.98 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.072 min
Delta R.T.: 0.000 min
Response: 377326236
Conc: 19.09 ng/ml

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168672BS	SDG No.:	Q2458
Lab Sample ID:	PB168672BS	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	100 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089275.D	1	07/01/25 08:30	07/01/25 16:23	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	18.7		0.13	1.70	ug/kg
319-85-7	beta-BHC	18.0		0.18	1.70	ug/kg
319-86-8	delta-BHC	19.2		0.39	1.70	ug/kg
58-89-9	gamma-BHC (Lindane)	18.6		0.14	1.70	ug/kg
76-44-8	Heptachlor	18.2		0.12	1.70	ug/kg
309-00-2	Aldrin	18.7		0.12	1.70	ug/kg
1024-57-3	Heptachlor epoxide	19.3		0.19	1.70	ug/kg
959-98-8	Endosulfan I	18.4		0.14	1.70	ug/kg
60-57-1	Dieldrin	18.4		0.14	1.70	ug/kg
72-55-9	4,4-DDE	18.4		0.14	1.70	ug/kg
72-20-8	Endrin	16.5		0.14	1.70	ug/kg
33213-65-9	Endosulfan II	17.9		0.29	1.70	ug/kg
72-54-8	4,4-DDD	18.7		0.15	1.70	ug/kg
1031-07-8	Endosulfan Sulfate	17.7		0.13	1.70	ug/kg
50-29-3	4,4-DDT	16.3		0.14	1.70	ug/kg
72-43-5	Methoxychlor	16.6		0.37	1.70	ug/kg
53494-70-5	Endrin ketone	18.3		0.19	1.70	ug/kg
7421-93-4	Endrin aldehyde	17.8		0.37	1.70	ug/kg
5103-71-9	alpha-Chlordane	18.2		0.12	1.70	ug/kg
5103-74-2	gamma-Chlordane	18.5		0.15	1.70	ug/kg
8001-35-2	Toxaphene	5.40	U	5.40	33.0	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	20.6		20 - 144	103%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.8		19 - 148	104%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168672BS	SDG No.:	Q2458
Lab Sample ID:	PB168672BS	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	100
Sample Wt/Vol:	30.03	Units:	g
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	Injection Volume :	
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089275.D	1	07/01/25 08:30	07/01/25 16:23	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089275.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 16:23
 Operator : AR\AJ
 Sample : PB168672BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB168672BS

Manual Integrations
 APPROVED

Reviewed By : Abdul Mirza 07/02/2025
 Supervised By : mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:47:27 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.556	2.880	52242928	351.9E6	18.311	20.757
28)	SA Decachlor...	9.081	8.074	80959011	385.8E6	20.613	19.517
Target Compounds							
2)	A alpha-BHC	4.006	3.393	326.9E6	1497.1E6	56.083	55.869
3)	MA gamma-BHC...	4.337	3.728	311.6E6	1371.6E6	55.828	55.401m
4)	MA Heptachlor	4.936	4.083	297.3E6	1322.0E6	54.531	52.798
5)	MB Aldrin	5.278	4.369	297.8E6	1337.7E6	56.057	55.236
6)	B beta-BHC	4.522	4.025	117.2E6	581.0E6	53.090	54.068
7)	B delta-BHC	4.771	4.262	295.7E6	1365.2E6	57.604	54.884
8)	B Heptachlo...	5.698	4.873	260.4E6	1269.3E6	53.983	57.966
9)	A Endosulfan I	6.081	5.248	247.1E6	1154.4E6	54.825	55.389
10)	B gamma-Chl...	5.953	5.125	265.6E6	1282.5E6	55.470	53.611m
11)	B alpha-Chl...	6.034	5.191	263.0E6	1248.2E6	54.364	54.512
12)	B 4,4'-DDE	6.203	5.376	241.3E6	1235.2E6	55.328	53.903
13)	MA Dieldrin	6.354	5.514	264.9E6	1254.1E6	55.218	54.038
14)	MA Endrin	6.581	5.790	204.4E6	1074.3E6	49.547	49.474
15)	B Endosulfa...	6.793	6.082	217.3E6	1077.5E6	53.823	53.350
16)	A 4,4'-DDD	6.712	5.930	193.2E6	1029.6E6	56.282	53.822
17)	MA 4,4'-DDT	7.029	6.185	184.9E6	970.4E6	48.798	47.851
18)	B Endrin al...	6.922	6.260	163.6E6	807.9E6	53.379	52.768
19)	B Endosulfa...	7.156	6.483	204.3E6	1036.6E6	53.237	52.832
20)	A Methoxychlor	7.500	6.755	98676499	537.4E6	48.835	49.906
21)	B Endrin ke...	7.637	6.993	223.4E6	1156.1E6	54.858	53.518
22)	Mirex	8.120	7.186	126.0E6	685.2E6	41.454	41.086

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089275.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 16:23
Operator : AR\AJ
Sample : PB168672BS
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PB168672BS

Manual Integrations

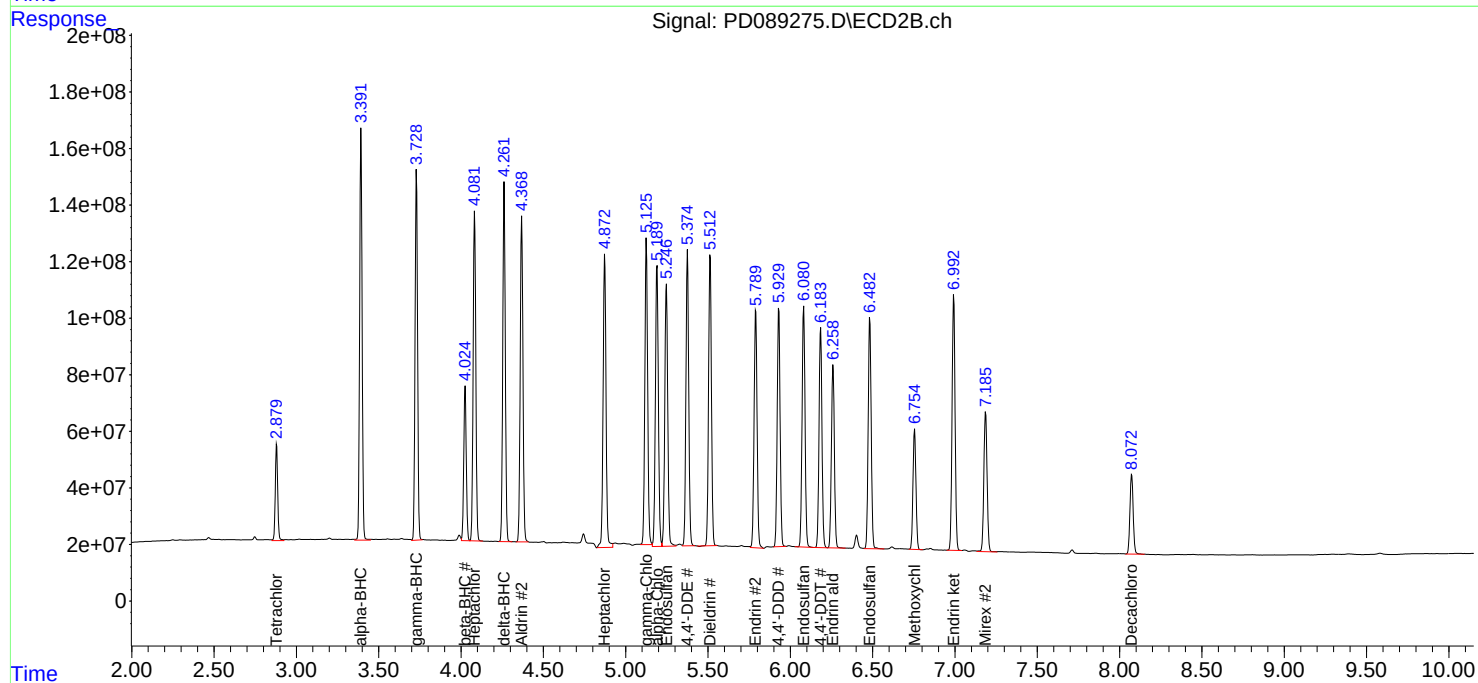
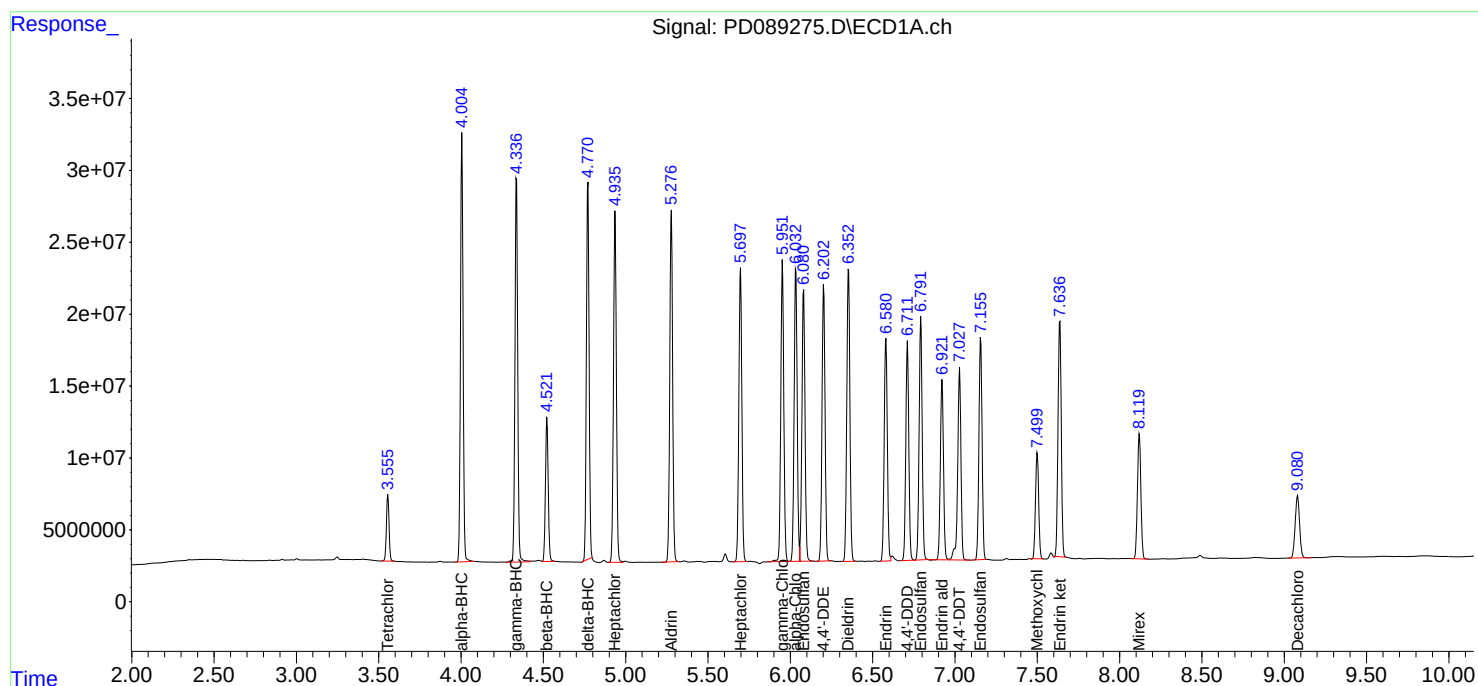
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:47:27 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BS	SDG No.:	Q2458
Lab Sample ID:	PB168718BS	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089330.D	1	07/03/25 08:58	07/03/25 14:09	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.56		0.0039	0.050	ug/L
319-85-7	beta-BHC	0.54		0.0049	0.050	ug/L
319-86-8	delta-BHC	0.59		0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.56		0.0037	0.050	ug/L
76-44-8	Heptachlor	0.54		0.0027	0.050	ug/L
309-00-2	Aldrin	0.56		0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.55		0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.56		0.0031	0.050	ug/L
60-57-1	Dieldrin	0.57		0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.56		0.0037	0.050	ug/L
72-20-8	Endrin	0.49		0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.57		0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.59		0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.55		0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.48		0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.44		0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.56		0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.56		0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.56		0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.57		0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	18.9		57 - 171	95%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.7		61 - 148	103%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BS	SDG No.:	Q2458
Lab Sample ID:	PB168718BS	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089330.D	1	07/03/25 08:58	07/03/25 14:09	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089330.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 14:09
 Operator : AR\AJ
 Sample : PB168718BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB168718BS

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/07/2025
 Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:17:40 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.551	2.881	50538235	350.2E6	17.713	20.659
28)	SA Decachlor...	9.076	8.071	73089221	373.6E6	18.609	18.897
Target Compounds							
2)	A alpha-BHC	4.000	3.393	323.5E6	1501.9E6	55.505	56.047
3)	MA gamma-BHC...	4.332	3.728	307.7E6	1377.9E6	55.121	55.657m
4)	MA Heptachlor	4.930	4.082	292.5E6	1312.4E6	53.651	52.416
5)	MB Aldrin	5.272	4.368	299.0E6	1344.8E6	56.285	55.532
6)	B beta-BHC	4.517	4.025	116.5E6	583.0E6	52.766	54.252
7)	B delta-BHC	4.766	4.262	300.8E6	1372.5E6	58.596	55.179
8)	B Heptachlo...	5.692	4.872	266.4E6	1215.1E6	55.226	55.487
9)	A Endosulfan I	6.076	5.247	251.8E6	1154.7E6	55.870	55.401
10)	B gamma-Chl...	5.948	5.125	273.4E6	1306.9E6	57.109	54.631
11)	B alpha-Chl...	6.029	5.190	268.9E6	1248.2E6	55.574	54.513
12)	B 4,4'-DDE	6.198	5.375	245.1E6	1256.6E6	56.211	54.834
13)	MA Dieldrin	6.349	5.512	271.3E6	1281.1E6	56.558	55.200
14)	MA Endrin	6.576	5.789	201.8E6	1032.9E6	48.903	47.570
15)	B Endosulfa...	6.788	6.080	228.7E6	1108.7E6	56.661	54.895
16)	A 4,4'-DDD	6.707	5.929	202.6E6	1072.7E6	59.018	56.077
17)	MA 4,4'-DDT	7.023	6.183	181.6E6	960.2E6	47.949	47.350
18)	B Endrin al...	6.917	6.258	171.8E6	843.6E6	56.069	55.099
19)	B Endosulfa...	7.151	6.482	211.0E6	1072.1E6	54.990	54.639
20)	A Methoxychlor	7.495	6.753	84312283	473.7E6	41.726	43.995
21)	B Endrin ke...	7.632	6.991	226.2E6	1192.7E6	55.534	55.212
22)	Mirex	8.116	7.185	136.5E6	752.8E6	44.916	45.134

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089330.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 14:09
Operator : AR\AJ
Sample : PB168718BS
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PB168718BS

Manual Integrations

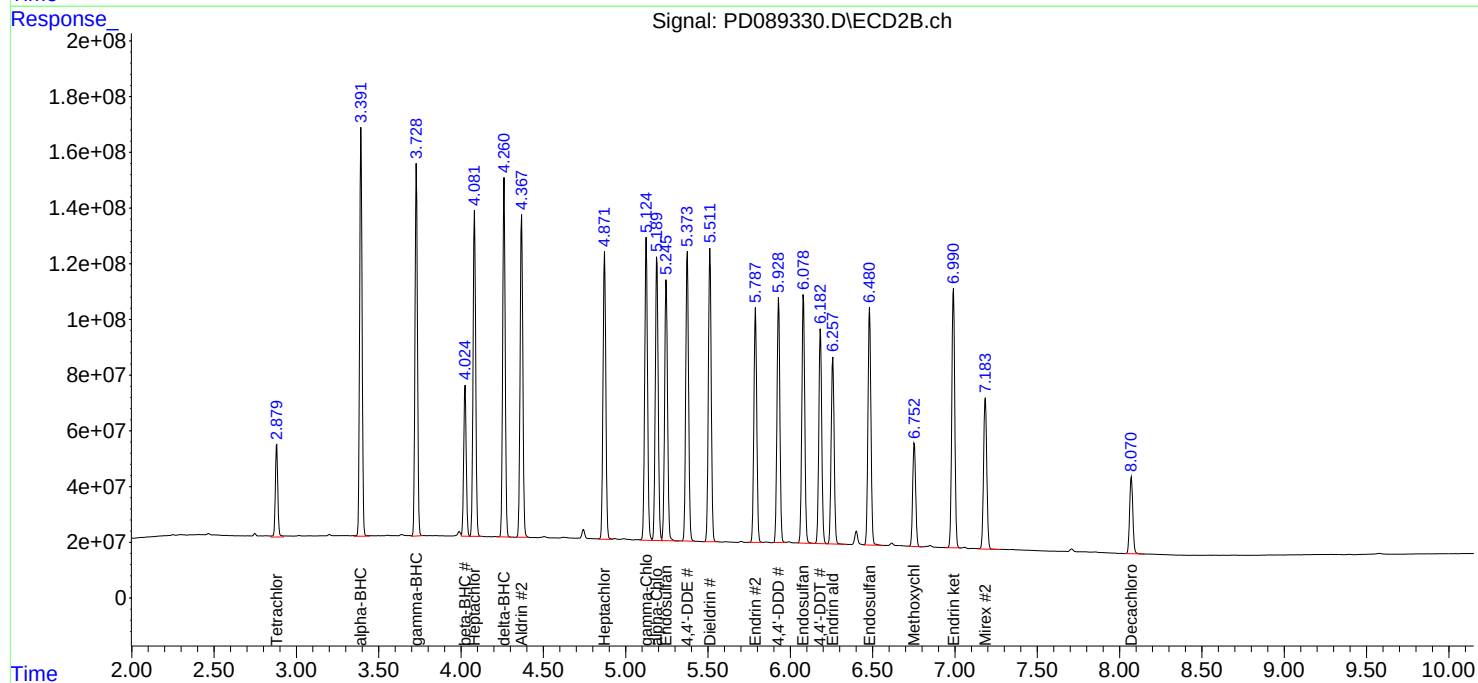
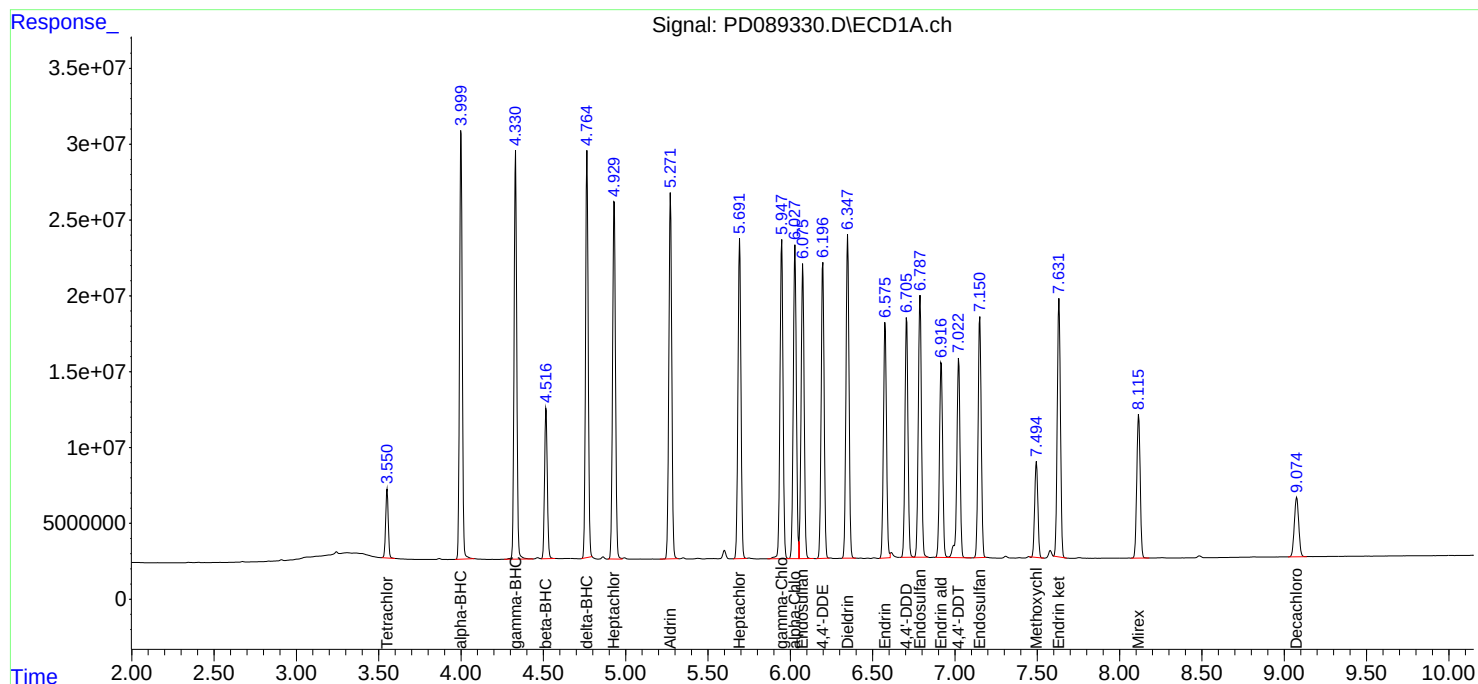
APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:17:40 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BSD	SDG No.:	Q2458
Lab Sample ID:	PB168718BSD	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089331.D	1	07/03/25 08:58	07/03/25 14:27	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
319-84-6	alpha-BHC	0.56		0.0039	0.050	ug/L
319-85-7	beta-BHC	0.54		0.0049	0.050	ug/L
319-86-8	delta-BHC	0.57		0.011	0.050	ug/L
58-89-9	gamma-BHC (Lindane)	0.55		0.0037	0.050	ug/L
76-44-8	Heptachlor	0.53		0.0027	0.050	ug/L
309-00-2	Aldrin	0.56		0.0036	0.050	ug/L
1024-57-3	Heptachlor epoxide	0.55		0.0096	0.050	ug/L
959-98-8	Endosulfan I	0.55		0.0031	0.050	ug/L
60-57-1	Dieldrin	0.56		0.0036	0.050	ug/L
72-55-9	4,4-DDE	0.55		0.0037	0.050	ug/L
72-20-8	Endrin	0.48		0.0032	0.050	ug/L
33213-65-9	Endosulfan II	0.56		0.0079	0.050	ug/L
72-54-8	4,4-DDD	0.58		0.0071	0.050	ug/L
1031-07-8	Endosulfan Sulfate	0.55		0.0037	0.050	ug/L
50-29-3	4,4-DDT	0.47		0.0035	0.050	ug/L
72-43-5	Methoxychlor	0.44		0.011	0.050	ug/L
53494-70-5	Endrin ketone	0.55		0.0093	0.050	ug/L
7421-93-4	Endrin aldehyde	0.55		0.011	0.050	ug/L
5103-71-9	alpha-Chlordane	0.55		0.0035	0.050	ug/L
5103-74-2	gamma-Chlordane	0.56		0.0039	0.050	ug/L
8001-35-2	Toxaphene	0.17	U	0.17	1.00	ug/L
SURROGATES						
2051-24-3	Decachlorobiphenyl	19.0		57 - 171	95%	SPK: 20
877-09-8	Tetrachloro-m-xylene	20.7		61 - 148	104%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	
Project:	South River WM Replacement	Date Received:	
Client Sample ID:	PB168718BSD	SDG No.:	Q2458
Lab Sample ID:	PB168718BSD	Matrix:	WATER
Analytical Method:	8081B	% Solid:	0 Decanted:
Sample Wt/Vol:	1000 Units: mL	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	3510C		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089331.D	1	07/03/25 08:58	07/03/25 14:27	PB168718

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
 Data File : PD089331.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 03 Jul 2025 14:27
 Operator : AR\AJ
 Sample : PB168718BSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB168718BSD

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/07/2025
 Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 04 02:17:56 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.555	2.880	50160508	351.7E6	17.581	20.745
28)	SA Decachlor...	9.081	8.074	72406437	375.1E6	18.435	18.976
Target Compounds							
2)	A alpha-BHC	4.005	3.393	321.4E6	1493.3E6	55.141	55.727
3)	MA gamma-BHC...	4.336	3.728	305.0E6	1368.9E6	54.649	55.296m
4)	MA Heptachlor	4.935	4.083	287.9E6	1303.7E6	52.815	52.068
5)	MB Aldrin	5.277	4.369	295.6E6	1338.5E6	55.634	55.271
6)	B beta-BHC	4.522	4.026	115.3E6	577.6E6	52.248	53.748
7)	B delta-BHC	4.770	4.263	293.0E6	1365.1E6	57.084	54.883
8)	B Heptachlo...	5.697	4.873	262.7E6	1210.5E6	54.451	55.277
9)	A Endosulfan I	6.081	5.248	248.7E6	1148.6E6	55.165	55.109
10)	B gamma-Chl...	5.952	5.126	269.9E6	1304.8E6	56.371	54.544
11)	B alpha-Chl...	6.033	5.191	265.9E6	1241.3E6	54.946	54.213
12)	B 4,4'-DDE	6.202	5.376	240.9E6	1250.5E6	55.233	54.569
13)	MA Dieldrin	6.353	5.514	268.4E6	1272.6E6	55.948	54.832
14)	MA Endrin	6.580	5.790	199.4E6	1039.4E6	48.330	47.866
15)	B Endosulfa...	6.792	6.082	226.0E6	1100.5E6	55.980	54.489
16)	A 4,4'-DDD	6.711	5.931	200.1E6	1070.2E6	58.301	55.945
17)	MA 4,4'-DDT	7.027	6.184	179.8E6	956.7E6	47.458	47.176
18)	B Endrin al...	6.922	6.258	170.0E6	842.7E6	55.493	55.036m
19)	B Endosulfa...	7.156	6.483	209.2E6	1071.3E6	54.528	54.601
20)	A Methoxychlor	7.499	6.755	83991927	473.6E6	41.567	43.977
21)	B Endrin ke...	7.636	6.992	223.8E6	1194.2E6	54.936	55.282
22)	Mirex	8.120	7.187	135.5E6	750.8E6	44.555	45.014

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070425\
Data File : PD089331.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 03 Jul 2025 14:27
Operator : AR\AJ
Sample : PB168718BSD
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PB168718BSD

Manual Integrations

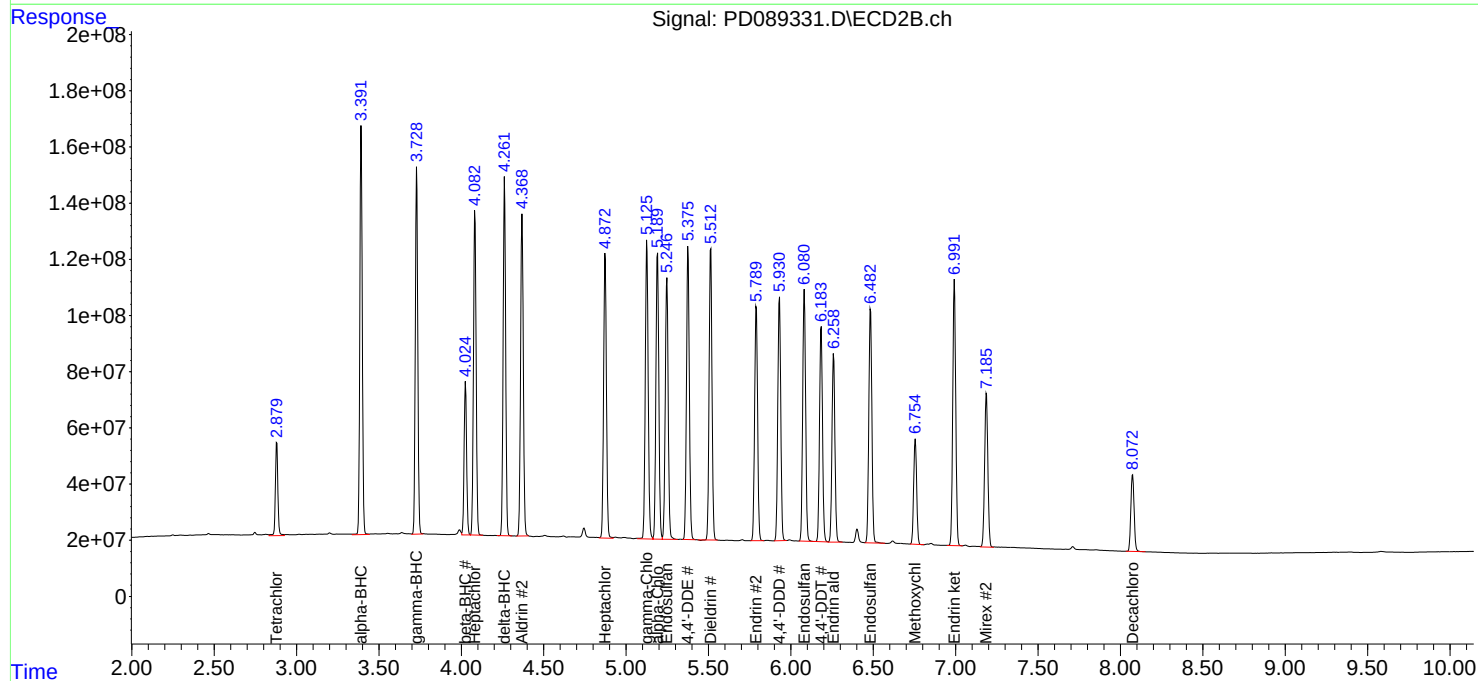
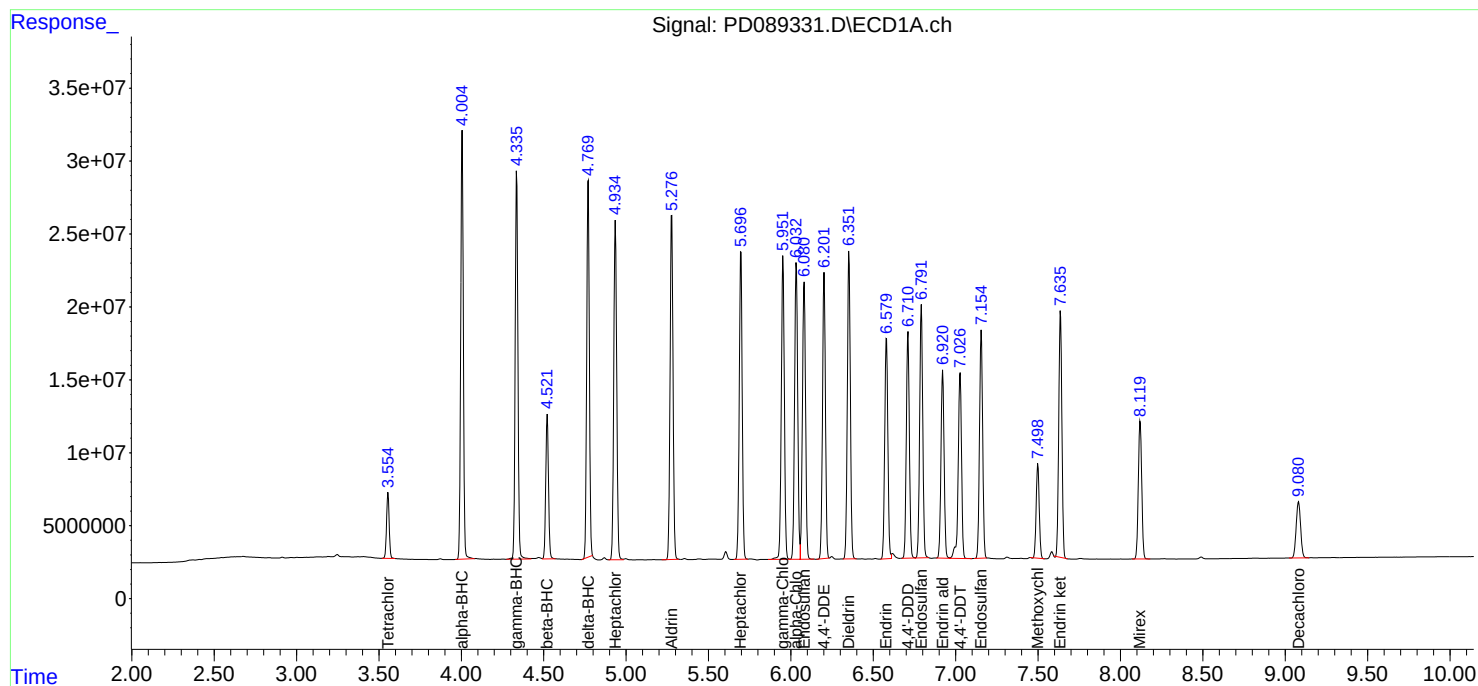
APPROVED

Reviewed By :Abdul Mirza 07/07/2025

Supervised By :mohammad ahmed 07/09/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 04 02:17:56 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67MS	SDG No.:	Q2458
Lab Sample ID:	Q2458-04MS	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7 Decanted:
Sample Wt/Vol:	30.03 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089291.D	1	07/01/25 08:30	07/01/25 20:29	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	17.7		0.14	1.90	ug/kg
319-85-7	beta-BHC	17.6		0.20	1.90	ug/kg
319-86-8	delta-BHC	17.9		0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	17.5		0.16	1.90	ug/kg
76-44-8	Heptachlor	16.5		0.13	1.90	ug/kg
309-00-2	Aldrin	18.0		0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	18.2		0.21	1.90	ug/kg
959-98-8	Endosulfan I	17.8		0.16	1.90	ug/kg
60-57-1	Dieldrin	17.7		0.16	1.90	ug/kg
72-55-9	4,4-DDE	18.0		0.16	1.90	ug/kg
72-20-8	Endrin	16.9		0.16	1.90	ug/kg
33213-65-9	Endosulfan II	17.2		0.32	1.90	ug/kg
72-54-8	4,4-DDD	18.4		0.17	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	16.4		0.14	1.90	ug/kg
50-29-3	4,4-DDT	14.2		0.16	1.90	ug/kg
72-43-5	Methoxychlor	13.4		0.41	1.90	ug/kg
53494-70-5	Endrin ketone	17.2		0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	16.8		0.41	1.90	ug/kg
5103-71-9	alpha-Chlordane	18.0		0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	18.3		0.17	1.90	ug/kg
8001-35-2	Toxaphene	6.00	U	6.00	36.8	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	16.5		20 - 144	82%	SPK: 20
877-09-8	Tetrachloro-m-xylene	21.3		19 - 148	107%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67MS	SDG No.:	Q2458
Lab Sample ID:	Q2458-04MS	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7
Sample Wt/Vol:	30.03	Units:	g
Soil Aliquot Vol:			uL
Extraction Type:		Test:	Pesticide-TCL
GPC Factor :	1.0	Injection Volume :	
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089291.D	1	07/01/25 08:30	07/01/25 20:29	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected
 LOQ = Limit of Quantitation
 MDL = Method Detection Limit
 LOD = Limit of Detection
 E = Value Exceeds Calibration Range
 P = Indicates >25% difference for detected concentrations between the two GC columns
 Q = indicates LCS control criteria did not meet requirements
 M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value
 B = Analyte Found in Associated Method Blank
 N = Presumptive Evidence of a Compound
 * = Values outside of QC limits
 D = Dilution
 S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.
 () = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089291.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 20:29
 Operator : AR\AJ
 Sample : Q2458-04MS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-67MS

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:51:36 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.880	55732238	361.4E6	19.534	21.320m
28)	SA Decachlor...	9.072	8.071	64068843	326.0E6	16.312	16.489
Target Compounds							
2)	A alpha-BHC	3.999	3.392	278.1E6	1280.4E6	47.705	47.780m
3)	MA gamma-BHC...	4.330	3.728	262.7E6	1164.8E6	47.062	47.051m
4)	MA Heptachlor	4.929	4.083	242.1E6	1097.5E6	44.409	43.834
5)	MB Aldrin	5.271	4.368	257.6E6	1170.8E6	48.478	48.344
6)	B beta-BHC	4.516	4.025	102.6E6	510.5E6	46.476	47.502
7)	B delta-BHC	4.764	4.262	248.2E6	1132.1E6	48.351	45.515
8)	B Heptachlo...	5.691	4.872	228.5E6	1073.2E6	47.363	49.007
9)	A Endosulfan I	6.074	5.247	215.9E6	988.2E6	47.891	47.414
10)	B gamma-Chl...	5.944	5.126	234.5E6	1176.0E6	48.972m	49.159
11)	B alpha-Chl...	6.027	5.190	235.0E6	1101.1E6	48.571	48.090
12)	B 4,4'-DDE	6.196	5.375	211.8E6	1104.5E6	48.566	48.199
13)	MA Dieldrin	6.347	5.513	227.9E6	1108.4E6	47.512	47.757
14)	MA Endrin	6.574	5.789	185.0E6	991.1E6	44.844	45.643
15)	B Endosulfa...	6.786	6.080	186.4E6	938.2E6	46.182	46.455
16)	A 4,4'-DDD	6.705	5.929	170.5E6	873.8E6	49.676	45.676
17)	MA 4,4'-DDT	7.021	6.183	143.1E6	776.5E6	37.788	38.291
18)	B Endrin al...	6.915	6.258	138.4E6	686.5E6	45.179	44.835
19)	B Endosulfa...	7.149	6.482	169.7E6	851.1E6	44.234	43.379
20)	A Methoxychlor	7.491	6.753	72698023	387.1E6	35.978m	35.951
21)	B Endrin ke...	7.630	6.991	188.9E6	950.6E6	46.374	44.004
22)	Mirex	8.114	7.184	135.9E6	728.3E6	44.705	43.669

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089291.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 20:29
Operator : AR\AJ
Sample : Q2458-04MS
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

TP-67MS

Manual Integrations

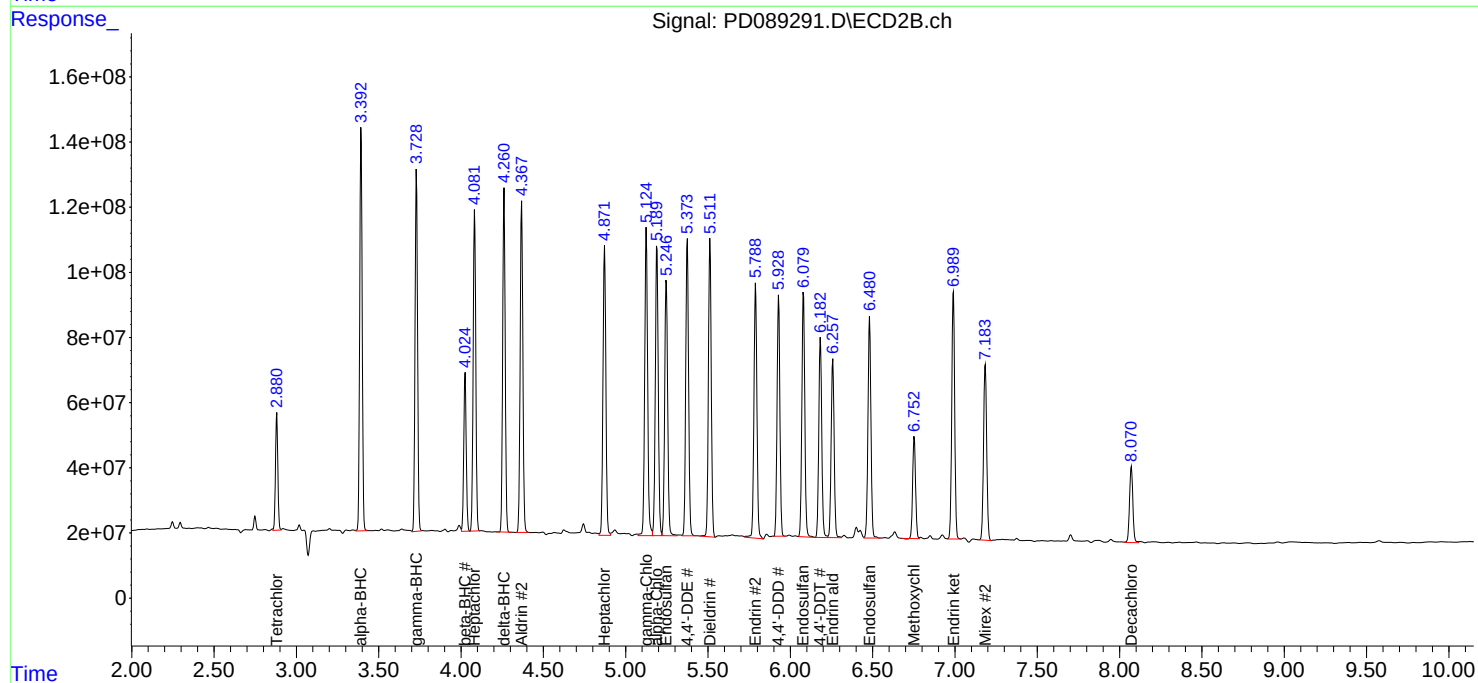
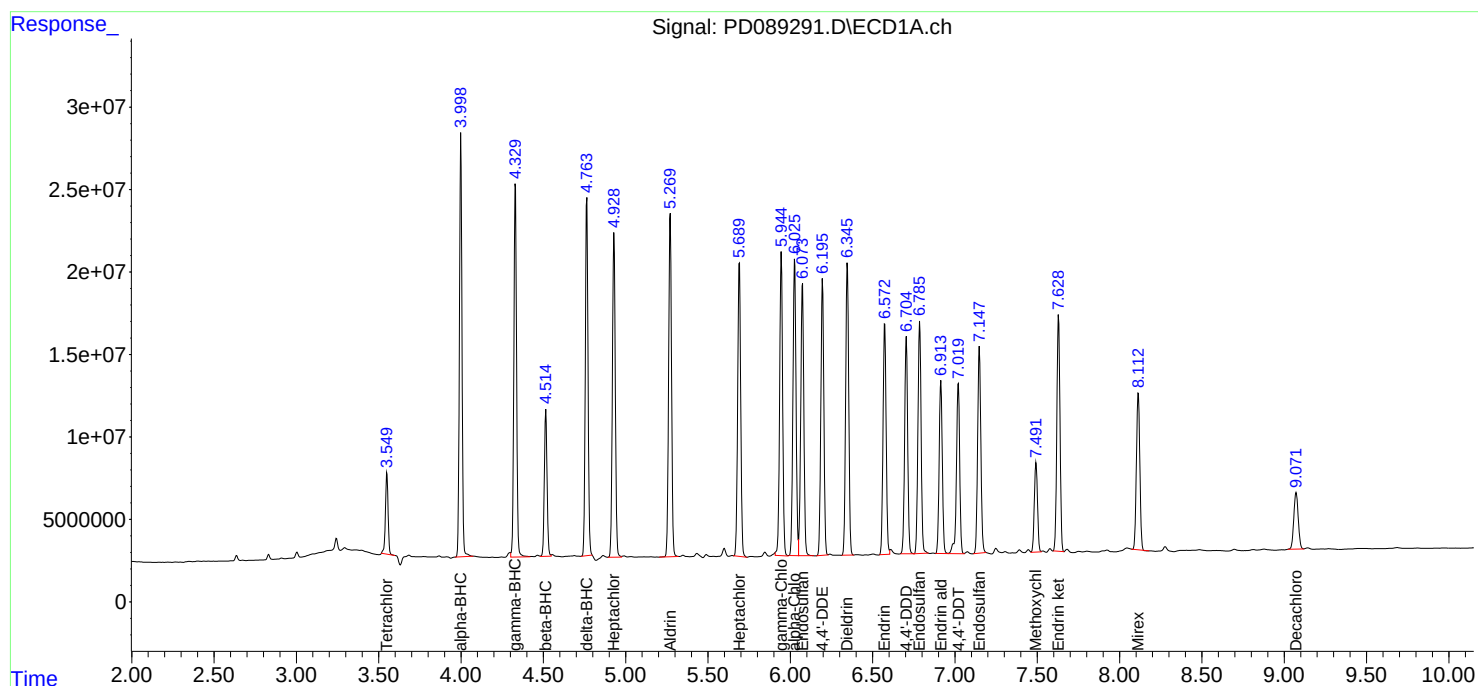
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:51:36 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67MSD	SDG No.:	Q2458
Lab Sample ID:	Q2458-04MSD	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7 Decanted:
Sample Wt/Vol:	30.07 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089292.D	1	07/01/25 08:30	07/01/25 20:43	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units(Dry Weight)
TARGETS						
319-84-6	alpha-BHC	18.2		0.14	1.90	ug/kg
319-85-7	beta-BHC	17.9		0.20	1.90	ug/kg
319-86-8	delta-BHC	19.0		0.43	1.90	ug/kg
58-89-9	gamma-BHC (Lindane)	18.2		0.16	1.90	ug/kg
76-44-8	Heptachlor	16.9		0.13	1.90	ug/kg
309-00-2	Aldrin	18.2		0.13	1.90	ug/kg
1024-57-3	Heptachlor epoxide	18.3		0.21	1.90	ug/kg
959-98-8	Endosulfan I	17.9		0.16	1.90	ug/kg
60-57-1	Dieldrin	17.9		0.16	1.90	ug/kg
72-55-9	4,4-DDE	18.0		0.16	1.90	ug/kg
72-20-8	Endrin	17.2		0.16	1.90	ug/kg
33213-65-9	Endosulfan II	17.4		0.32	1.90	ug/kg
72-54-8	4,4-DDD	19.0		0.17	1.90	ug/kg
1031-07-8	Endosulfan Sulfate	17.1		0.14	1.90	ug/kg
50-29-3	4,4-DDT	15.0		0.16	1.90	ug/kg
72-43-5	Methoxychlor	14.2		0.41	1.90	ug/kg
53494-70-5	Endrin ketone	17.7		0.21	1.90	ug/kg
7421-93-4	Endrin aldehyde	17.3		0.41	1.90	ug/kg
5103-71-9	alpha-Chlordane	18.2		0.13	1.90	ug/kg
5103-74-2	gamma-Chlordane	18.4		0.17	1.90	ug/kg
8001-35-2	Toxaphene	6.00	U	6.00	36.7	ug/kg
SURROGATES						
2051-24-3	Decachlorobiphenyl	16.6		20 - 144	83%	SPK: 20
877-09-8	Tetrachloro-m-xylene	21.6		19 - 148	108%	SPK: 20

Report of Analysis

Client:	CDM Smith	Date Collected:	06/27/25
Project:	South River WM Replacement	Date Received:	06/27/25
Client Sample ID:	TP-67MSD	SDG No.:	Q2458
Lab Sample ID:	Q2458-04MSD	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	89.7 Decanted:
Sample Wt/Vol:	30.07 Units: g	Final Vol:	10000 uL
Soil Aliquot Vol:	uL	Test:	Pesticide-TCL
Extraction Type:		Injection Volume :	
GPC Factor :	1.0 PH :		
Prep Method :	SW3541B		

File ID/Qc Batch:	Dilution:	Prep Date	Date Analyzed	Prep Batch ID
PD089292.D	1	07/01/25 08:30	07/01/25 20:43	PB168672

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
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Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

P = Indicates >25% difference for detected concentrations between the two GC columns

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

* = Values outside of QC limits

D = Dilution

S = Indicates estimated value where valid five-point calibration was not performed prior to analyte detection in sample.

() = Laboratory InHouse Limit

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
 Data File : PD089292.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 01 Jul 2025 20:43
 Operator : AR\AJ
 Sample : Q2458-04MSD
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TP-67MSD

Manual Integrations
 APPROVED

Reviewed By :Abdul Mirza 07/02/2025
 Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Jul 02 01:51:48 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
 Quant Title : GC Extractables
 QLast Update : Wed Jun 18 05:39:05 2025
 Response via : Initial Calibration
 Integrator: ChemStation

Volume Inj. : 1 µl
 Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
 Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

	Compound	RT#1	RT#2	Resp#1	Resp#2	ng/ml	ng/ml

System Monitoring Compounds							
1)	SA Tetrachlo...	3.550	2.880	56098549	367.0E6	19.662	21.650m
28)	SA Decachlor...	9.073	8.072	63612232	327.6E6	16.196	16.573
Target Compounds							
2)	A alpha-BHC	4.000	3.392	285.8E6	1310.3E6	49.030	48.897m
3)	MA gamma-BHC...	4.331	3.728	274.7E6	1202.7E6	49.222	48.579m
4)	MA Heptachlor	4.929	4.083	248.9E6	1131.5E6	45.647	45.192
5)	MB Aldrin	5.271	4.369	260.3E6	1180.6E6	48.993	48.751
6)	B beta-BHC	4.516	4.026	103.6E6	519.0E6	46.932	48.296
7)	B delta-BHC	4.764	4.262	263.1E6	1197.3E6	51.252	48.136
8)	B Heptachlo...	5.691	4.872	230.8E6	1080.9E6	47.854	49.359
9)	A Endosulfan I	6.074	5.247	217.9E6	1003.8E6	48.353	48.161
10)	B gamma-Chl...	5.946	5.125	234.7E6	1189.6E6	49.023	49.725
11)	B alpha-Chl...	6.027	5.190	237.5E6	1114.3E6	49.078	48.664
12)	B 4,4'-DDE	6.196	5.375	211.6E6	1105.2E6	48.531	48.230
13)	MA Dieldrin	6.347	5.511	230.9E6	1118.5E6	48.120	48.195m
14)	MA Endrin	6.574	5.789	187.9E6	1007.1E6	45.550	46.381
15)	B Endosulfa...	6.786	6.080	188.9E6	949.8E6	46.792	47.027
16)	A 4,4'-DDD	6.705	5.929	176.0E6	899.1E6	51.278	47.000
17)	MA 4,4'-DDT	7.021	6.183	150.3E6	820.2E6	39.689	40.445
18)	B Endrin al...	6.915	6.258	142.7E6	709.2E6	46.571	46.319
19)	B Endosulfa...	7.150	6.482	176.6E6	896.8E6	46.018	45.706
20)	A Methoxychlor	7.493	6.752	75092398	412.4E6	37.163	38.301m
21)	B Endrin ke...	7.630	6.991	194.5E6	990.0E6	47.754	45.830
22)	Mirex	8.114	7.184	137.4E6	736.9E6	45.209	44.184

(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD070225\
Data File : PD089292.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 01 Jul 2025 20:43
Operator : AR\AJ
Sample : Q2458-04MSD
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

TP-67MSD

Manual Integrations

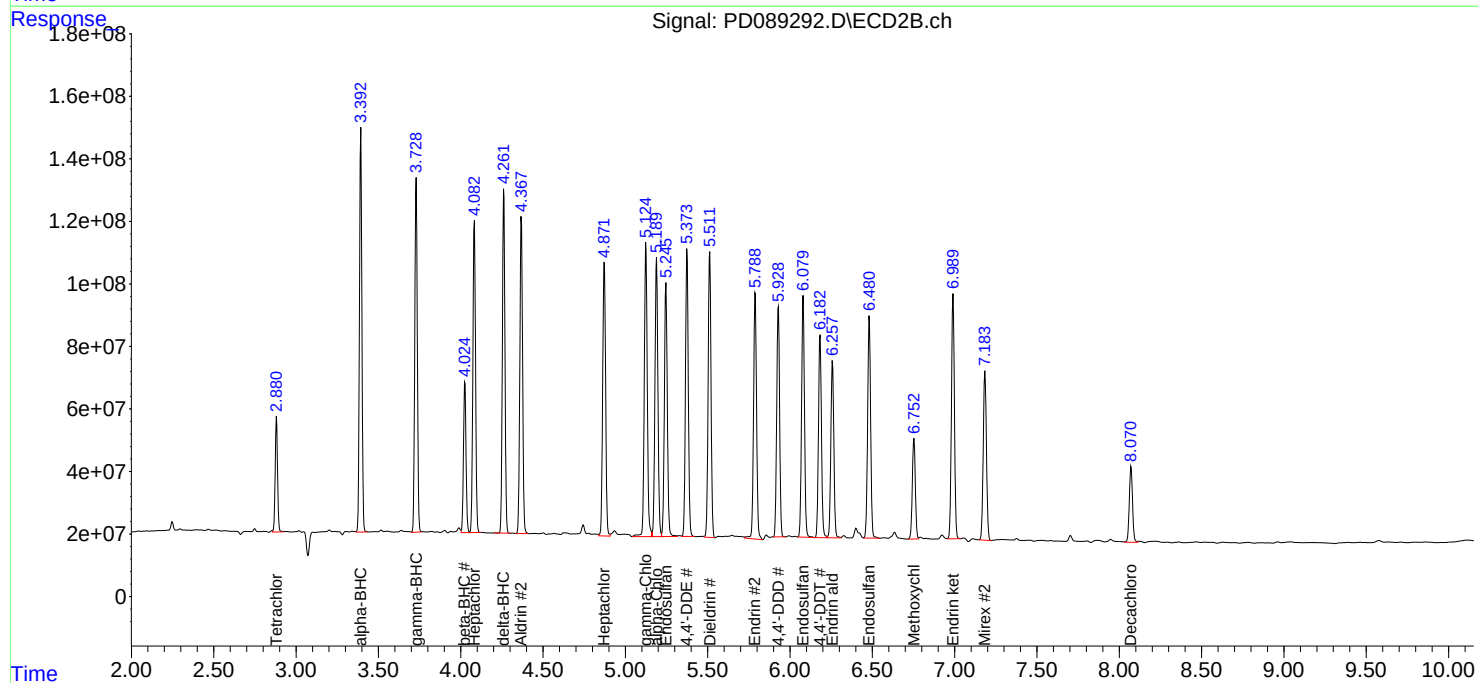
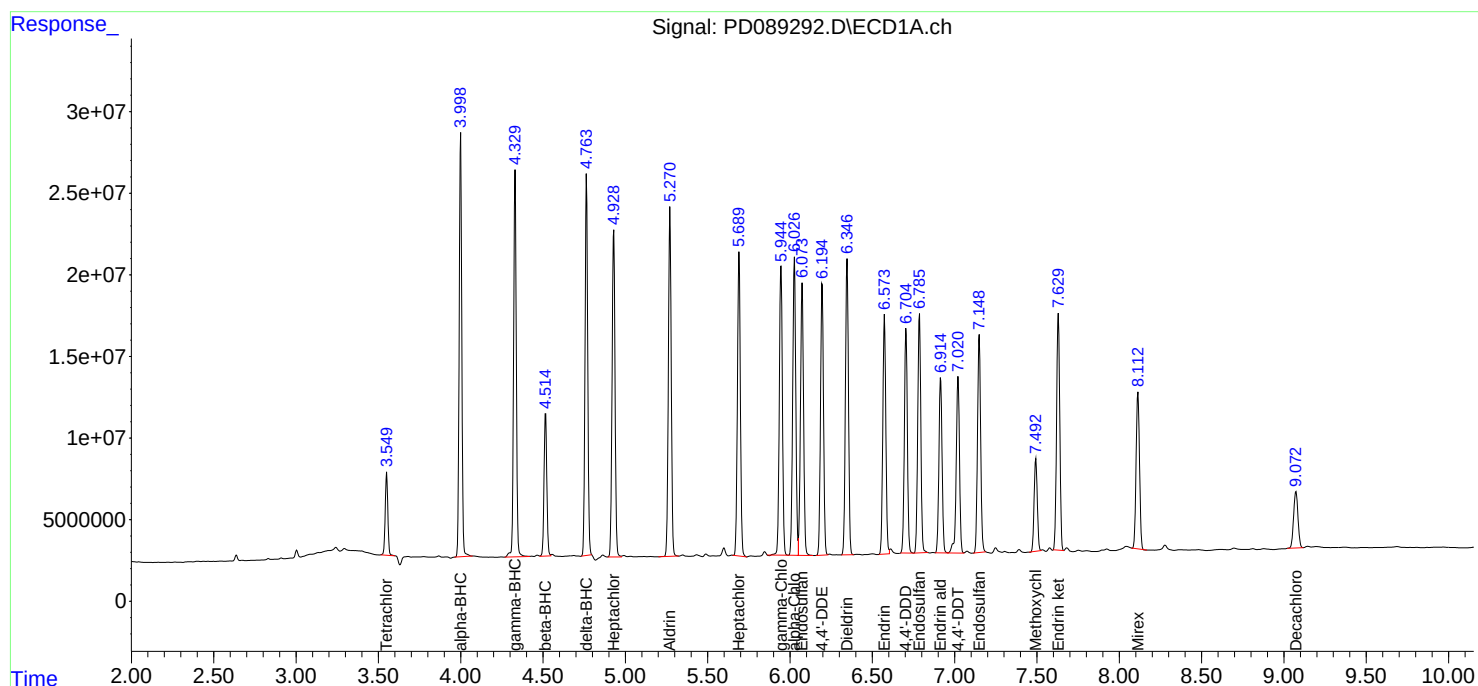
APPROVED

Reviewed By :Abdul Mirza 07/02/2025

Supervised By :mohammad ahmed 07/04/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Jul 02 01:51:48 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD061825.M
Quant Title : GC Extractables
QLast Update : Wed Jun 18 05:39:05 2025
Response via : Initial Calibration
Integrator: ChemStation

Volume Inj. : 1 µl
Signal #1 Phase : ZB-MR1 Signal #2 Phase: ZB-MR2
Signal #1 Info : 30M x 0.32mm x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Manual Integration Report

Sequence:	PD061825	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PD088991.D	4,4"-DDD	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	4,4"-DDD #2	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	beta-BHC	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	Endrin aldehyde	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	Endrin aldehyde #2	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	Endrin ketone	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PEM	PD088991.D	Endrin ketone #2	Abdul	6/18/2025 8:40:15 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PSTDICC025	PD088996.D	Endrin #2	Abdul	6/18/2025 8:40:24 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PSTDICC005	PD088997.D	Endrin #2	Abdul	6/18/2025 8:40:30 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software
PSTDICC005	PD088997.D	Endrin ketone #2	Abdul	6/18/2025 8:40:30 AM	mohammad	6/19/2025 2:43:55	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PD089265.D	4,4"-DDD	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089265.D	4,4"-DDE	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089265.D	4,4"-DDE #2	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089265.D	alpha-BHC #2	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089265.D	Endrin	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089265.D	gamma-BHC (Lindane)	Abdul	7/2/2025 8:57:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	delta-BHC	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	Dieldrin #2	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	Endrin aldehyde	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	gamma-Chlordane #2	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	Heptachlor epoxide #2	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089266.D	Mirex #2	Abdul	7/2/2025 8:57:53 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PB168672BS	PD089275.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:58:16 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PB168672BS	PD089275.D	gamma-Chlordane #2	Abdul	7/2/2025 8:58:16 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	alpha-Chlordane #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	Dieldrin #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	Endosulfan I #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	gamma-Chlordane #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089277.D	Heptachlor epoxide #2	Abdul	7/2/2025 8:58:19 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-01	PD089284.D	Decachlorobiphenyl	Abdul	7/2/2025 8:58:33 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-01	PD089284.D	Dieldrin	Abdul	7/2/2025 8:58:33 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-01	PD089284.D	Dieldrin #2	Abdul	7/2/2025 8:58:33 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-01	PD089284.D	Endrin #2	Abdul	7/2/2025 8:58:33 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-01	PD089284.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:58:33 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
Q2458-03	PD089286.D	Decachlorobiphenyl	Abdul	7/2/2025 8:58:36 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-03	PD089286.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:58:36 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089288.D	4,4"-DDE #2	Abdul	7/2/2025 8:58:39 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089288.D	Endrin	Abdul	7/2/2025 8:58:39 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PEM	PD089288.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:58:39 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089289.D	Dieldrin #2	Abdul	7/2/2025 8:58:43 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089289.D	gamma-Chlordane #2	Abdul	7/2/2025 8:58:43 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089289.D	Heptachlor epoxide #2	Abdul	7/2/2025 8:58:43 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04	PD089290.D	4,4"-DDE #2	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04	PD089290.D	4,4"-DDT	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04	PD089290.D	alpha-Chlordane #2	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04	PD089290.D	gamma-Chlordane	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04	PD089290.D	gamma-Chlordane #2	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
Q2458-04	PD089290.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:58:46 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MS	PD089291.D	alpha-BHC #2	Abdul	7/2/2025 8:58:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MS	PD089291.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:58:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MS	PD089291.D	gamma-Chlordane	Abdul	7/2/2025 8:58:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MS	PD089291.D	Methoxychlor	Abdul	7/2/2025 8:58:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MS	PD089291.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:58:50 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MSD	PD089292.D	alpha-BHC #2	Abdul	7/2/2025 8:58:54 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MSD	PD089292.D	Dieldrin #2	Abdul	7/2/2025 8:58:54 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MSD	PD089292.D	gamma-BHC (Lindane) #2	Abdul	7/2/2025 8:58:54 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MSD	PD089292.D	Methoxychlor #2	Abdul	7/2/2025 8:58:54 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-04MSD	PD089292.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:58:54 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-05	PD089293.D	4,4"-DDT	Abdul	7/2/2025 3:42:44 PM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-05	PD089293.D	4,4"-DDT #2	Abdul	7/2/2025 3:42:44 PM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
Q2458-06	PD089294.D	4,4"-DDD	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	4,4"-DDD #2	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	4,4"-DDT	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	alpha-Chlordane	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	alpha-Chlordane #2	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	gamma-Chlordane	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-06	PD089294.D	gamma-Chlordane #2	Abdul	7/2/2025 8:59:05 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-07	PD089295.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:59:11 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-08	PD089296.D	4,4"-DDD	Abdul	7/2/2025 8:59:15 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-08	PD089296.D	alpha-Chlordane #2	Abdul	7/2/2025 8:59:15 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-08	PD089296.D	gamma-Chlordane	Abdul	7/2/2025 8:59:15 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-08	PD089296.D	gamma-Chlordane #2	Abdul	7/2/2025 8:59:15 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
Q2458-08	PD089296.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:59:15 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:	pd070225	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
Q2458-09	PD089297.D	Tetrachloro-m-xylene #2	Abdul	7/2/2025 8:59:20 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software
PSTDCCC050	PD089299.D	Dieldrin #2	Abdul	7/2/2025 8:59:24 AM	mohammad	7/4/2025 4:31:21	Peak Integrated by Software

Manual Integration Report

Sequence:

pd070425

Instrument

ECD_d

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
I.BLK	PD089326.D	Tetrachloro-m-xylene #2	Abdul	7/7/2025 8:15:12 AM	 	 	Peak Integrated by Software
PEM	PD089327.D	4,4"-DDD	Abdul	7/7/2025 8:14:38 AM	 	 	Peak Integrated by Software
PEM	PD089327.D	4,4"-DDD #2	Abdul	7/7/2025 8:14:38 AM	 	 	Peak Integrated by Software
PEM	PD089327.D	4,4"-DDE	Abdul	7/7/2025 8:14:38 AM	 	 	Peak Integrated by Software
PEM	PD089327.D	4,4"-DDE #2	Abdul	7/7/2025 8:14:38 AM	 	 	Peak Integrated by Software
PSTDCCC050	PD089328.D	Dieldrin #2	Abdul	7/7/2025 8:13:26 AM	 	 	Peak Integrated by Software
PSTDCCC050	PD089328.D	gamma-BHC (Lindane) #2	Abdul	7/7/2025 8:13:26 AM	 	 	Peak Integrated by Software
PSTDCCC050	PD089328.D	gamma-Chlordane #2	Abdul	7/7/2025 8:13:26 AM	 	 	Peak Integrated by Software
PB168718BL	PD089329.D	Tetrachloro-m-xylene #2	Abdul	7/7/2025 8:16:43 AM	 	 	Peak Integrated by Software
PB168718BS	PD089330.D	gamma-BHC (Lindane) #2	Abdul	7/7/2025 8:16:33 AM	 	 	Peak Integrated by Software
PB168718BSD	PD089331.D	Endrin aldehyde #2	Abdul	7/7/2025 8:16:09 AM	 	 	Peak Integrated by Software
PB168718BSD	PD089331.D	gamma-BHC (Lindane) #2	Abdul	7/7/2025 8:16:09 AM	 	 	Peak Integrated by Software
PSTDCCC050	PD089338.D	Dieldrin #2	Abdul	7/7/2025 8:14:07 AM	 	 	Peak Integrated by Software



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900, Fax : 908 789 8922

Manual Integration Report

Sequence:

pd070425

Instrument

ECD_d

Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PSTDCCC050	PD089338.D	gamma-BHC (Lindane) #2	Abdul	7/7/2025 8:14:07 AM	 	 	Peak Integrated by Software

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD061825

Review By	Abdul	Review On	6/18/2025 8:40:56 AM
Supervise By	mohammad	Supervise On	6/19/2025 2:43:55 AM
SubDirectory	PD061825	HP Acquire Method	HP Processing Method pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,PP24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	Sampleld	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD088989.D	17 Jun 2025 14:58	AR\AJ	Ok
2	I.BLK	PD088990.D	17 Jun 2025 15:11	AR\AJ	Ok
3	PEM	PD088991.D	17 Jun 2025 15:25	AR\AJ	Ok,M
4	RESCHK	PD088992.D	17 Jun 2025 15:39	AR\AJ	Ok
5	PSTDICC100	PD088993.D	17 Jun 2025 15:52	AR\AJ	Ok
6	PSTDICC075	PD088994.D	17 Jun 2025 16:06	AR\AJ	Ok
7	PSTDICC050	PD088995.D	17 Jun 2025 16:20	AR\AJ	Ok
8	PSTDICC025	PD088996.D	17 Jun 2025 16:33	AR\AJ	Ok,M
9	PSTDICC005	PD088997.D	17 Jun 2025 16:47	AR\AJ	Ok,M
10	PCHLORICC1000	PD088998.D	17 Jun 2025 17:00	AR\AJ	Ok
11	PCHLORICC750	PD088999.D	17 Jun 2025 17:14	AR\AJ	Ok
12	PCHLORICC500	PD089000.D	17 Jun 2025 17:28	AR\AJ	Ok
13	PCHLORICC250	PD089001.D	17 Jun 2025 17:41	AR\AJ	Ok
14	PCHLORICC050	PD089002.D	17 Jun 2025 17:55	AR\AJ	Ok
15	PTOXICC1000	PD089003.D	17 Jun 2025 18:09	AR\AJ	Ok
16	PTOXICC750	PD089004.D	17 Jun 2025 18:22	AR\AJ	Ok
17	PTOXICC500	PD089005.D	17 Jun 2025 18:36	AR\AJ	Ok
18	PTOXICC250	PD089006.D	17 Jun 2025 18:50	AR\AJ	Ok
19	PTOXICC100	PD089007.D	17 Jun 2025 19:03	AR\AJ	Ok,M
20	PSTDICV050	PD089008.D	17 Jun 2025 19:17	AR\AJ	Ok
21	PCHLORICV500	PD089009.D	17 Jun 2025 19:31	AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD061825

Review By	Abdul	Review On	6/18/2025 8:40:56 AM
Supervise By	mohammad	Supervise On	6/19/2025 2:43:55 AM
SubDirectory	PD061825	HP Acquire Method	HP Processing Method pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,PP24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

22	PTOXICV500	PD089010.D	17 Jun 2025 19:44	AR\AJ	Ok
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M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070225

Review By	Abdul	Review On	7/2/2025 9:00:09 AM
Supervise By	mohammad	Supervise On	7/4/2025 4:31:21 AM
SubDirectory	PD070225	HP Acquire Method	HP Processing Method pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,PP24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD089263.D	01 Jul 2025 11:46	AR\AJ	Ok
2	I.BLK	PD089264.D	01 Jul 2025 11:59	AR\AJ	Ok
3	PEM	PD089265.D	01 Jul 2025 12:13	AR\AJ	Ok,M
4	PSTDCCC050	PD089266.D	01 Jul 2025 13:35	AR\AJ	Ok,M
5	Q2447-04	PD089267.D	01 Jul 2025 13:52	AR\AJ	Ok,M
6	Q2447-06	PD089268.D	01 Jul 2025 14:06	AR\AJ	Ok
7	PB168655BL	PD089269.D	01 Jul 2025 14:19	AR\AJ	Ok,M
8	PB168672BL	PD089270.D	01 Jul 2025 14:33	AR\AJ	Ok
9	PB168672BS	PD089271.D	01 Jul 2025 14:47	AR\AJ	Not Ok
10	Q2464-01	PD089272.D	01 Jul 2025 15:00	AR\AJ	Ok,M
11	Q2459-01	PD089273.D	01 Jul 2025 15:14	AR\AJ	Ok,M
12	Q2469-01	PD089274.D	01 Jul 2025 15:28	AR\AJ	Ok,M
13	PB168672BS	PD089275.D	01 Jul 2025 16:23	AR\AJ	Ok,M
14	I.BLK	PD089276.D	01 Jul 2025 16:37	AR\AJ	Ok
15	PSTDCCC050	PD089277.D	01 Jul 2025 16:50	AR\AJ	Ok,M
16	PB168676BL	PD089278.D	01 Jul 2025 17:04	AR\AJ	Ok
17	PB168676BS	PD089279.D	01 Jul 2025 17:18	AR\AJ	Ok,M
18	PB168571TB	PD089280.D	01 Jul 2025 17:32	AR\AJ	Ok
19	Q2409-02	PD089281.D	01 Jul 2025 17:45	AR\AJ	Ok
20	Q2409-02MS	PD089282.D	01 Jul 2025 17:59	AR\AJ	Ok,M
21	Q2409-02MSD	PD089283.D	01 Jul 2025 18:13	AR\AJ	Ok,M

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070225

Review By	Abdul	Review On	7/2/2025 9:00:09 AM
Supervise By	mohammad	Supervise On	7/4/2025 4:31:21 AM
SubDirectory	PD070225	HP Acquire Method	HP Processing Method pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,PP24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

22	Q2458-01	PD089284.D	01 Jul 2025 18:26	AR\AJ	Ok,M
23	Q2458-02	PD089285.D	01 Jul 2025 18:40	AR\AJ	Ok
24	Q2458-03	PD089286.D	01 Jul 2025 18:53	AR\AJ	Ok,M
25	I.BLK	PD089287.D	01 Jul 2025 19:07	AR\AJ	Ok
26	PEM	PD089288.D	01 Jul 2025 19:21	AR\AJ	Ok,M
27	PSTDCCC050	PD089289.D	01 Jul 2025 19:34	AR\AJ	Ok,M
28	Q2458-04	PD089290.D	01 Jul 2025 20:15	AR\AJ	Ok,M
29	Q2458-04MS	PD089291.D	01 Jul 2025 20:29	AR\AJ	Ok,M
30	Q2458-04MSD	PD089292.D	01 Jul 2025 20:43	AR\AJ	Ok,M
31	Q2458-05	PD089293.D	01 Jul 2025 20:57	AR\AJ	Ok,M
32	Q2458-06	PD089294.D	01 Jul 2025 21:10	AR\AJ	Ok,M
33	Q2458-07	PD089295.D	01 Jul 2025 21:24	AR\AJ	Ok,M
34	Q2458-08	PD089296.D	01 Jul 2025 21:37	AR\AJ	Ok,M
35	Q2458-09	PD089297.D	01 Jul 2025 21:51	AR\AJ	Ok,M
36	I.BLK	PD089298.D	01 Jul 2025 22:05	AR\AJ	Ok
37	PSTDCCC050	PD089299.D	01 Jul 2025 22:59	AR\AJ	Ok,M

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070425

Review By	Review On
Supervise By	Supervise On
SubDirectory PD070425	HP Acquire Method HP Processing Method pd061825 8081
STD. NAME	STD REF.#
Tune/Reschk	PP24433,PP24651
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,PP24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284
CCC	PP24261,PP24273,PP24279,PP24284
Internal Standard/PEM	
ICV/I.BLK	PP24273,PP24279,PP24284
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD089325.D	03 Jul 2025 09:05	AR\AJ	Ok
2	I.BLK	PD089326.D	03 Jul 2025 09:19	AR\AJ	Ok,NS
3	PEM	PD089327.D	03 Jul 2025 09:32	AR\AJ	Ok,NS
4	PSTDCCC050	PD089328.D	03 Jul 2025 09:46	AR\AJ	Ok,NS
5	PB168718BL	PD089329.D	03 Jul 2025 13:55	AR\AJ	Ok,NS
6	PB168718BS	PD089330.D	03 Jul 2025 14:09	AR\AJ	Ok,NS
7	PB168718BSD	PD089331.D	03 Jul 2025 14:27	AR\AJ	Ok,NS
8	Q2458-10	PD089332.D	03 Jul 2025 14:40	AR\AJ	Ok
9	PB168725BL	PD089333.D	03 Jul 2025 14:54	AR\AJ	Ok
10	PB168725BS	PD089334.D	03 Jul 2025 15:07	AR\AJ	Ok,NS
11	Q2487-09	PD089335.D	03 Jul 2025 15:21	AR\AJ	Ok
12	Q2487-10	PD089336.D	03 Jul 2025 15:35	AR\AJ	Ok,NS
13	I.BLK	PD089337.D	03 Jul 2025 15:48	AR\AJ	Ok
14	PSTDCCC050	PD089338.D	03 Jul 2025 16:02	AR\AJ	Ok,NS

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD061825

Review By	Abdul	Review On	6/18/2025 8:40:56 AM
Supervise By	mohammad	Supervise On	6/19/2025 2:43:55 AM
SubDirectory	PD061825	HP Acquire Method	HP Processing Method pd061825 8081

STD. NAME	STD REF.#
Tune/Reschk	PP24433,PP24651
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284
CCC	PP24261,PP24273,PP24279,PP24284
Internal Standard/PEM	
ICV/I.BLK	PP24273,PP24279,PP24284
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	Sampleld	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD088989.D	17 Jun 2025 14:58		AR\AJ	Ok
2	I.BLK	I.BLK	PD088990.D	17 Jun 2025 15:11		AR\AJ	Ok
3	PEM	PEM	PD088991.D	17 Jun 2025 15:25		AR\AJ	Ok,M
4	RESCHK	RESCHK	PD088992.D	17 Jun 2025 15:39		AR\AJ	Ok
5	PSTDICC100	PSTDICC100	PD088993.D	17 Jun 2025 15:52		AR\AJ	Ok
6	PSTDICC075	PSTDICC075	PD088994.D	17 Jun 2025 16:06		AR\AJ	Ok
7	PSTDICC050	PSTDICC050	PD088995.D	17 Jun 2025 16:20		AR\AJ	Ok
8	PSTDICC025	PSTDICC025	PD088996.D	17 Jun 2025 16:33		AR\AJ	Ok,M
9	PSTDICC005	PSTDICC005	PD088997.D	17 Jun 2025 16:47		AR\AJ	Ok,M
10	PCHLORICC1000	PCHLORICC1000	PD088998.D	17 Jun 2025 17:00		AR\AJ	Ok
11	PCHLORICC750	PCHLORICC750	PD088999.D	17 Jun 2025 17:14		AR\AJ	Ok
12	PCHLORICC500	PCHLORICC500	PD089000.D	17 Jun 2025 17:28		AR\AJ	Ok
13	PCHLORICC250	PCHLORICC250	PD089001.D	17 Jun 2025 17:41		AR\AJ	Ok
14	PCHLORICC050	PCHLORICC050	PD089002.D	17 Jun 2025 17:55		AR\AJ	Ok
15	PTOXICC1000	PTOXICC1000	PD089003.D	17 Jun 2025 18:09		AR\AJ	Ok
16	PTOXICC750	PTOXICC750	PD089004.D	17 Jun 2025 18:22		AR\AJ	Ok
17	PTOXICC500	PTOXICC500	PD089005.D	17 Jun 2025 18:36		AR\AJ	Ok
18	PTOXICC250	PTOXICC250	PD089006.D	17 Jun 2025 18:50		AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD061825

Review By	Abdul	Review On	6/18/2025 8:40:56 AM
Supervise By	mohammad	Supervise On	6/19/2025 2:43:55 AM
SubDirectory	PD061825	HP Acquire Method	HP Processing Method
			pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

19	PTOXICC100	PTOXICC100	PD089007.D	17 Jun 2025 19:03		AR\AJ	Ok,M
20	PSTDICV050	ICVPD061825	PD089008.D	17 Jun 2025 19:17		AR\AJ	Ok
21	PCHLORICV500	ICVPD061825CHLOR	PD089009.D	17 Jun 2025 19:31		AR\AJ	Ok
22	PTOXICV500	ICVPD061825TOX	PD089010.D	17 Jun 2025 19:44		AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070225

Review By	Abdul	Review On	7/2/2025 9:00:09 AM
Supervise By	mohammad	Supervise On	7/4/2025 4:31:21 AM
SubDirectory	PD070225	HP Acquire Method	HP Processing Method pd061825 8081

STD. NAME	STD REF.#
Tune/Reschk	PP24433,PP24651
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284
CCC	PP24261,PP24273,PP24279,PP24284
Internal Standard/PEM	
ICV/I.BLK	PP24273,PP24279,PP24284
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	SampleId	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD089263.D	01 Jul 2025 11:46		AR\AJ	Ok
2	I.BLK	I.BLK	PD089264.D	01 Jul 2025 11:59		AR\AJ	Ok
3	PEM	PEM	PD089265.D	01 Jul 2025 12:13		AR\AJ	Ok,M
4	PSTDCCC050	PSTDCCC050	PD089266.D	01 Jul 2025 13:35		AR\AJ	Ok,M
5	Q2447-04	COMP-2	PD089267.D	01 Jul 2025 13:52	check cleanup	AR\AJ	Ok,M
6	Q2447-06	LAW-25-0100	PD089268.D	01 Jul 2025 14:06	check cleanup	AR\AJ	Ok
7	PB168655BL	PB168655BL	PD089269.D	01 Jul 2025 14:19		AR\AJ	Ok,M
8	PB168672BL	PB168672BL	PD089270.D	01 Jul 2025 14:33		AR\AJ	Ok
9	PB168672BS	PB168672BS	PD089271.D	01 Jul 2025 14:47	Surrogate and some compounds recovery fail	AR\AJ	Not Ok
10	Q2464-01	OR-3-063025	PD089272.D	01 Jul 2025 15:00		AR\AJ	Ok,M
11	Q2459-01	TP-1	PD089273.D	01 Jul 2025 15:14		AR\AJ	Ok,M
12	Q2469-01	WC-1	PD089274.D	01 Jul 2025 15:28		AR\AJ	Ok,M
13	PB168672BS	PB168672BS	PD089275.D	01 Jul 2025 16:23		AR\AJ	Ok,M
14	I.BLK	I.BLK	PD089276.D	01 Jul 2025 16:37		AR\AJ	Ok
15	PSTDCCC050	PSTDCCC050	PD089277.D	01 Jul 2025 16:50		AR\AJ	Ok,M
16	PB168676BL	PB168676BL	PD089278.D	01 Jul 2025 17:04		AR\AJ	Ok
17	PB168676BS	PB168676BS	PD089279.D	01 Jul 2025 17:18		AR\AJ	Ok,M

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070225

Review By	Abdul	Review On	7/2/2025 9:00:09 AM
Supervise By	mohammad	Supervise On	7/4/2025 4:31:21 AM
SubDirectory	PD070225	HP Acquire Method	HP Processing Method pd061825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24433,PP24651		
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284		
CCC	PP24261,PP24273,PP24279,PP24284		
Internal Standard/PEM			
ICV/I.BLK	PP24273,PP24279,PP24284		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

18	PB168571TB	PB168571TB	PD089280.D	01 Jul 2025 17:32		AR\AJ	Ok
19	Q2409-02	COP-SOIL-PILE	PD089281.D	01 Jul 2025 17:45		AR\AJ	Ok
20	Q2409-02MS	COP-SOIL-PILEMS	PD089282.D	01 Jul 2025 17:59		AR\AJ	Ok,M
21	Q2409-02MSD	COP-SOIL-PILEMSD	PD089283.D	01 Jul 2025 18:13		AR\AJ	Ok,M
22	Q2458-01	TP-76	PD089284.D	01 Jul 2025 18:26		AR\AJ	Ok,M
23	Q2458-02	TP-55	PD089285.D	01 Jul 2025 18:40		AR\AJ	Ok
24	Q2458-03	TP-68	PD089286.D	01 Jul 2025 18:53		AR\AJ	Ok,M
25	I.BLK	I.BLK	PD089287.D	01 Jul 2025 19:07		AR\AJ	Ok
26	PEM	PEM	PD089288.D	01 Jul 2025 19:21		AR\AJ	Ok,M
27	PSTDCCC050	PSTDCCC050	PD089289.D	01 Jul 2025 19:34		AR\AJ	Ok,M
28	Q2458-04	TP-67	PD089290.D	01 Jul 2025 20:15		AR\AJ	Ok,M
29	Q2458-04MS	TP-67MS	PD089291.D	01 Jul 2025 20:29		AR\AJ	Ok,M
30	Q2458-04MSD	TP-67MSD	PD089292.D	01 Jul 2025 20:43		AR\AJ	Ok,M
31	Q2458-05	TP-66	PD089293.D	01 Jul 2025 20:57		AR\AJ	Ok,M
32	Q2458-06	TP-60	PD089294.D	01 Jul 2025 21:10		AR\AJ	Ok,M
33	Q2458-07	TP-62	PD089295.D	01 Jul 2025 21:24		AR\AJ	Ok,M
34	Q2458-08	TP-63	PD089296.D	01 Jul 2025 21:37		AR\AJ	Ok,M
35	Q2458-09	TP-59	PD089297.D	01 Jul 2025 21:51		AR\AJ	Ok,M
36	I.BLK	I.BLK	PD089298.D	01 Jul 2025 22:05		AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070225

Review By	Abdul	Review On	7/2/2025 9:00:09 AM
Supervise By	mohammad	Supervise On	7/4/2025 4:31:21 AM
SubDirectory	PD070225	HP Acquire Method	HP Processing Method pd061825 8081

STD. NAME	STD REF.#
Tune/Reschk	PP24433,PP24651
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284
CCC	PP24261,PP24273,PP24279,PP24284
Internal Standard/PEM	
ICV/I.BLK	PP24273,PP24279,PP24284
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

37	PSTDCCC050	PSTDCCC050	PD089299.D	01 Jul 2025 22:59		ARIAJ	Ok,M
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M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD070425

Review By	Review On
Supervise By	Supervise On
SubDirectory	PD070425
HP Acquire Method	HP Processing Method
	pd061825 8081
STD. NAME	STD REF.#
Tune/Reschk	PP24433,PP24651
Initial Calibration Stds	PP24260,PP24261,PP24262,PP24269,PP24266,PP24267,PP24268,PP24269,PP24270,PP24271,PP24272,PP24273,PP24274,PP24275,PP24277,P P24278,PP24279,PP24280,PP24281,PP24282,PP24283,PP24284
CCC	PP24261,PP24273,PP24279,PP24284
Internal Standard/PEM	
ICV/I.BLK	PP24273,PP24279,PP24284
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	Sampleld	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD089325.D	03 Jul 2025 09:05		AR\AJ	Ok
2	I.BLK	I.BLK	PD089326.D	03 Jul 2025 09:19		AR\AJ	Ok,NS
3	PEM	PEM	PD089327.D	03 Jul 2025 09:32		AR\AJ	Ok,NS
4	PSTDCCC050	PSTDCCC050	PD089328.D	03 Jul 2025 09:46		AR\AJ	Ok,NS
5	PB168718BL	PB168718BL	PD089329.D	03 Jul 2025 13:55		AR\AJ	Ok,NS
6	PB168718BS	PB168718BS	PD089330.D	03 Jul 2025 14:09		AR\AJ	Ok,NS
7	PB168718BSD	PB168718BSD	PD089331.D	03 Jul 2025 14:27		AR\AJ	Ok,NS
8	Q2458-10	FB-06272025	PD089332.D	03 Jul 2025 14:40		AR\AJ	Ok
9	PB168725BL	PB168725BL	PD089333.D	03 Jul 2025 14:54		AR\AJ	Ok
10	PB168725BS	PB168725BS	PD089334.D	03 Jul 2025 15:07		AR\AJ	Ok,NS
11	Q2487-09	G4(0-6)	PD089335.D	03 Jul 2025 15:21		AR\AJ	Ok
12	Q2487-10	G4(6-12)	PD089336.D	03 Jul 2025 15:35		AR\AJ	Ok,NS
13	I.BLK	I.BLK	PD089337.D	03 Jul 2025 15:48		AR\AJ	Ok
14	PSTDCCC050	PSTDCCC050	PD089338.D	03 Jul 2025 16:02		AR\AJ	Ok,NS

M : Manual Integration

PERCENT SOLID

Supervisor: Iwona
Analyst: jignesh
Date: 7/1/2025

OVENTEMP IN Celsius(°C): 107
Time IN: 17:25
In Date: 06/30/2025
Weight Check 1.0g: 1.00
Weight Check 10g: 10.00
OvenID: M OVEN#1

OVENTEMP OUT Celsius(°C): 104
Time OUT: 08:22
Out Date: 07/01/2025
Weight Check 1.0g: 1.00
Weight Check 10g: 10.00
BalanceID: M SC-4
Thermometer ID: % SOLID-OVEN

QC:LB136327

Lab ID	Client SampleID	Dish #	Dish Wt(g) (A)	Sample Wt(g)	Dish + Sample Wt(g) (B)	Dish+Dry Sample Wt(g) (C)	% Solid	Comments
Q2451-01	Y2310-0409-1-1	1	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-02	Y2310-0409-1-2	2	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-03	Y2307-0279-1-1	3	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-04	Y2307-0279-1-2	4	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-05	KQA080Z-1-1	5	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-06	KQA080Z-1-2	6	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-07	HIA-925Q-1-1	7	1.00	1.00	2.00	2.00	100.0	PILC
Q2451-08	HIA-925Q-1-2	8	1.00	1.00	2.00	2.00	100.0	PILC
Q2458-01	TP-76	9	1.15	10.84	11.99	10.97	90.6	
Q2458-02	TP-55	10	1.18	10.49	11.67	10.77	91.4	
Q2458-03	TP-68	11	1.18	10.68	11.86	11.04	92.3	
Q2458-04	TP-67	12	1.19	10.16	11.35	10.3	89.7	
Q2458-05	TP-66	13	1.19	10.18	11.37	10.18	88.3	
Q2458-06	TP-60	14	1.15	10.83	11.98	11.17	92.5	
Q2458-07	TP-62	15	1.15	10.84	11.99	11.02	91.1	
Q2458-08	TP-63	16	1.19	10.82	12.01	10.54	86.4	
Q2458-09	TP-59	17	1.15	10.70	11.85	9.35	76.6	
Q2459-01	TP-1	18	1.19	10.74	11.93	10.66	88.2	
Q2459-02	TP-1-EPH	19	1.15	10.88	12.03	10.8	88.7	
Q2459-03	TP-1-VOC	20	1.16	10.83	11.99	10.9	89.9	
Q2459-04	TP-1	21	1.19	10.74	11.93	10.66	88.2	
Q2462-02	60425-A	22	1.15	10.31	11.46	9.96	85.5	
Q2462-03	60425-B	23	1.14	10.66	11.8	7.91	63.5	
Q2462-04	60425-AB	24	1.19	10.57	11.76	9.66	80.1	
Q2462-05	50725	25	1.00	1.00	2.00	2.00	100.0	debris
Q2462-06	61625-A	26	1.12	10.70	11.82	9.5	78.3	
Q2462-07	61625-B	27	1.18	10.37	11.55	8.29	68.6	
Q2462-08	61625-C	28	1.19	10.65	11.84	9.11	74.4	



PERCENT SOLID

Supervisor: Iwona
Analyst: jignesh
Date: 7/1/2025

OVENTEMP IN Celsius(°C): 107
Time IN: 17:25
In Date: 06/30/2025
Weight Check 1.0g: 1.00
Weight Check 10g: 10.00
OvenID: M OVEN#1

OVENTEMP OUT Celsius(°C): 104
Time OUT: 08:22
Out Date: 07/01/2025
Weight Check 1.0g: 1.00
Weight Check 10g: 10.00
BalanceID: M SC-4
Thermometer ID: % SOLID-OVEN

QC:LB136327

Lab ID	Client SampleID	Dish #	Dish Wt(g) (A)	Sample Wt(g)	Dish + Sample Wt(g) (B)	Dish+Dry Sample Wt(g) (C)	% Solid	Comments
Q2462-09	61625-ABC	29	1.19	10.56	11.75	9.26	76.4	
Q2462-10	62725	30	1.00	1.00	2.00	2.00	100.0	debris
Q2464-01	OR-3-063025	31	1.12	10.76	11.88	11.57	97.1	
Q2464-02	OR-3-063025-E2	32	1.18	10.50	11.68	11.3	96.4	
Q2466-01	61825	33	1.00	1.00	2.00	2.00	100.0	WIPE SAMPLE
Q2466-02	62625	34	1.00	1.00	2.00	2.00	100.0	WIPE SAMPLE
Q2467-01	1A-1B-1C-Fire Stopher Caulk-Red	35	1.00	1.00	2.00	2.00	100.0	Caulk sample
Q2469-01	WC-1	36	1.15	10.13	11.28	9.8	85.4	
Q2469-02	WC-1-EPH	37	1.19	10.36	11.55	10.14	86.4	
Q2469-03	WC-1-VOC	38	1.15	10.84	11.99	10.74	88.5	
Q2469-04	WC-1	39	1.15	10.13	11.28	9.8	85.4	

$$\% \text{ Solid} = \frac{(C-A) * 100}{(B-A)}$$

WORKLIST(Hardcopy Internal Chain)

136327

WorkList Name : %1-063025

WorkList ID : 190450

Department : Wet-Chemistry

Date : 06-30-2025 08:06:09

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q2451-01	Y2310-0409-1-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-02	Y2310-0409-1-2	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-03	Y2307-0279-1-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-04	Y2307-0279-1-2	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-05	KQA08OZ-1-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-06	KQA08OZ-1-2	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-07	HIA-925Q-1-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2451-08	HIA-925Q-1-2	Solid	Percent Solids	Cool 4 deg C	PSEG03	D51	06/27/2025	Chemtech -SO
Q2458-01	TP-76	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/26/2025	Chemtech -SO
Q2458-02	TP-55	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-03	TP-68	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-04	TP-67	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-05	TP-66	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-06	TP-60	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-07	TP-62	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-08	TP-63	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2458-09	TP-59	Solid	Percent Solids	Cool 4 deg C	CAMP02	D51	06/27/2025	Chemtech -SO
Q2459-01	TP-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	A51	06/28/2025	Chemtech -SO
Q2459-02	TP-1-EPH	Solid	Percent Solids	Cool 4 deg C	PSEG03	A51	06/28/2025	Chemtech -SO
Q2459-03	TP-1-VOC	Solid	Percent Solids	Cool 4 deg C	PSEG03	A51	06/28/2025	Chemtech -SO
Q2459-04	TP-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	A51	06/28/2025	Chemtech -SO

Date/Time 06/30/25 16:00

Raw Sample Received by: JWC

Raw Sample Relinquished by: JWC

Date/Time 06/30/25

Raw Sample Received by: JWC

Raw Sample Relinquished by: JWC

WORKLIST(Hardcopy Internal Chain)

130327

WorkList Name : %1-063025 WorkList ID : 190450 Department : Wet-Chemistry Date : 06-30-2025 08:06:09

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q2462-02	60425-A	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-03	60425-3	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-04	60425-AB	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-05	50725	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-06	61625-A	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-07	61625-B	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-08	61625-C	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-09	60425-ABC	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2462-10	62725	Solid	Percent Solids	Cool 4 deg C	PSEG03	A12	06/30/2025	Chemtech -SO
Q2464-01	OR-3-063025	Solid	Percent Solids	Cool 4 deg C	PSEG05	A21	06/30/2025	Chemtech -SO
Q2464-02	OR-3-063025	Solid	Percent Solids	Cool 4 deg C	PSEG05	A21	06/30/2025	Chemtech -SO
Q2466-01	61825	Solid	Percent Solids	Cool 4 deg C	PSEG03	A11	06/30/2025	Chemtech -SO
Q2466-02	62625	Solid	Percent Solids	Cool 4 deg C	PSEG03	A11	06/30/2025	Chemtech -SO
Q2467-01	1A-1B-1C-Fire Stopper Caulk-R	Solid	Percent Solids	Cool 4 deg C	ATCE02	A61	06/26/2025	Chemtech -SO
Q2469-01	WC-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	A61	06/30/2025	Chemtech -SO
Q2469-02	WC-1-EPH	Solid	Percent Solids	Cool 4 deg C	PSEG03	A61	06/30/2025	Chemtech -SO
Q2469-03	WC-1-VOC	Solid	Percent Solids	Cool 4 deg C	PSEG03	A61	06/30/2025	Chemtech -SO
Q2469-04	WC-1	Solid	Percent Solids	Cool 4 deg C	PSEG03	A61	06/30/2025	Chemtech -SO

06/30/25 16:00

Date/Time Raw Sample Received by: SP CSM Raw Sample Relinquished by: SP CSM

SOP ID: M3541-ASE Extraction-14

Clean Up SOP #: Florisil

Extraction Start Date : 07/01/2025

Matrix : Solid

Extraction Start Time : 08:30

Weigh By: EH

Extraction By: RJ

Extraction End Date : 07/01/2025

Balance check: RJ

Filter By: RJ

Extraction End Time : 11:40

Balance ID: EX-SC-2

pH Meter ID: N/A

Concentration By: EH

pH Strip Lot#: N/A

Hood ID: 3,7

Supervisor By : RUPESH

Extraction Method: ☐ Separatory Funnel ☐ Continuous Liquid/Liquid ☐ Sonication ☐ Waste Dilution ☒ Soxhlet

Standard Name	MLS USED	Concentration ug/mL	STD REF. # FROM LOG
Spike Sol 1	1.0ML	500 PPB	PP24627
Surrogate	1.0ML	200 PPB	PP24663
N/A	N/A	N/A	N/A
N/A	N/A	N/A	N/A
N/A	N/A	N/A	N/A

Chemical Used	ML/SAMPLE USED	Lot Number
Hexane/Acetone/1:1	N/A	EP2613
Baked Na2SO4	N/A	EP2624
Sand	N/A	E2865
Hexane	N/A	E3947
Florisil	N/A	E3927
9:1 Hexane:Acetone Mixture	N/A	EP2596
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A

Extraction Conformance/Non-Conformance Comments:

40ML Vial Lot # 03-40BTS723.

KD Bath ID: N/A

Envap ID: NE VAP-02

KD Bath Temperature: N/A

Envap Temperature: 40 °C

Date / Time	Prepped Sample Relinquished By/Location	Received By/Location
7/1/25	RS (Ext Lab)	Y-P-PCB+PCB
11:45	Preparation Group	Analysis Group

Analytical Method: M3541-ASE Extraction-14

Concentration Date: 07/01/2025

Sample ID	Client Sample ID	Test	(g)/ mL	PH	Surr/Spike By:		Final Vol. (mL)	JarID	Comments	Prep Pos
					AddedBy	VerifiedBy				
PB168672BL	PBLK672	Pesticide-TCL	30.01	N/A	ritesh	Evelyn	10			U1-1
PB168672BS	PLCS672	Pesticide-TCL	30.03	N/A	ritesh	Evelyn	10			2
Q2458-01	TP-76	Pesticide-TCL	30.04	N/A	ritesh	Evelyn	10	E		3
Q2458-02	TP-55	Pesticide-TCL	30.03	N/A	ritesh	Evelyn	10	E		4
Q2458-03	TP-68	Pesticide-TCL	30.01	N/A	ritesh	Evelyn	10	E		5
Q2458-04	TP-67	Pesticide-TCL	30.06	N/A	ritesh	Evelyn	10	E		6
Q2458-04MS	TP-67MS	Pesticide-TCL	30.03	N/A	ritesh	Evelyn	10	E		U2-1
Q2458-04MS D	TP-67MSD	Pesticide-TCL	30.07	N/A	ritesh	Evelyn	10	E		2
Q2458-05	TP-66	Pesticide-TCL	30.06	N/A	ritesh	Evelyn	10	E		3
Q2458-06	TP-60	Pesticide-TCL	30.05	N/A	ritesh	Evelyn	10	E		4
Q2458-07	TP-62	Pesticide-TCL	30.04	N/A	ritesh	Evelyn	10	E		5
Q2458-08	TP-63	Pesticide-TCL	30.08	N/A	ritesh	Evelyn	10	E		6
Q2458-09	TP-59	Pesticide-TCL	30.03	N/A	ritesh	Evelyn	10	E		U3-1
Q2459-01	TP-1	Pesticide-TCL	30.01	N/A	ritesh	Evelyn	10	D		2
Q2464-01	OR-3-063025	Pesticide-TCL	30.09	N/A	ritesh	Evelyn	10	E		3
Q2469-01	WC-1	Pesticide-TCL	30.07	N/A	ritesh	Evelyn	10	D		4

* Extracts relinquished on the same date as received.

RS
7/1/25

168672
2469891
8:30

WORKLIST(Hardcopy Internal Chain)

WorkList Name : Q2458 WorkList ID : 190471 Department : Extraction Date : 07-01-2025 08:24:37

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q2458-01	TP-76	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/26/2025	8081B
Q2458-02	TP-55	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/26/2025	8081B
Q2458-03	TP-68	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-04	TP-67	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-05	TP-66	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-06	TP-60	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-07	TP-62	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-08	TP-63	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2458-09	TP-59	Solid	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B
Q2459-01	TP-1	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	A51	06/28/2025	8081B
Q2464-01	OR-3-063025	Solid	Pesticide-TCL	Cool 4 deg C	PSEG05	A21	06/30/2025	8081B
Q2469-01	WC-1	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	A61	06/30/2025	8081B

Date/Time 07/01/25 8:25
Raw Sample Received by: RJ CFA-106J
Raw Sample Relinquished by: OP SN

Date/Time 07/01/25 8:50
Raw Sample Received by: OP SN
Raw Sample Relinquished by: RJ CFA-106J

SOP ID: M3510C,3580A-Extraction Pesticide-17

Clean Up SOP #: Florisil **Extraction Start Date :** 07/03/2025

Matrix : Water **Extraction Start Time :** 08:58

Weigh By: N/A **Extraction By:** RS **Extraction End Date :** 07/03/2025

Balance check: N/A **Filter By:** RS **Extraction End Time :** 13:30

Balance ID: N/A **pH Meter ID:** N/A **Concentration By:** EH

pH Strip Lot#: E3880 **Hood ID:** 4,5,6,7 **Supervisor By :** RUPESH

Extraction Method: ☒ Separatory Funnel ☐ Continuous Liquid/Liquid ☐ Sonication ☐ Waste Dilution ☐ Soxhlet

Standard Name	MLS USED	Concentration ug/mL	STD REF. # FROM LOG
Spike Sol 1	1.0ML	500 PPB	PP24627
Surrogate	1.0ML	200 PPB	PP24663
N/A	N/A	N/A	N/A
N/A	N/A	N/A	N/A
N/A	N/A	N/A	N/A

Chemical Used	ML/SAMPLE USED	Lot Number
Methylene Chloride	N/A	E3943
Baked Na2SO4	N/A	EP2624
Hexane	N/A	E3947
Florisil	N/A	E3927
9:1 Hexane:Acetone Mixture	N/A	EP2596
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A
N/A	N/A	N/A

Extraction Conformance/Non-Conformance Comments:

40 ML Vial lot# 03-40 BTS723.

KD Bath ID: WATER BATH-1,2 **Envap ID:** NEVAP-02

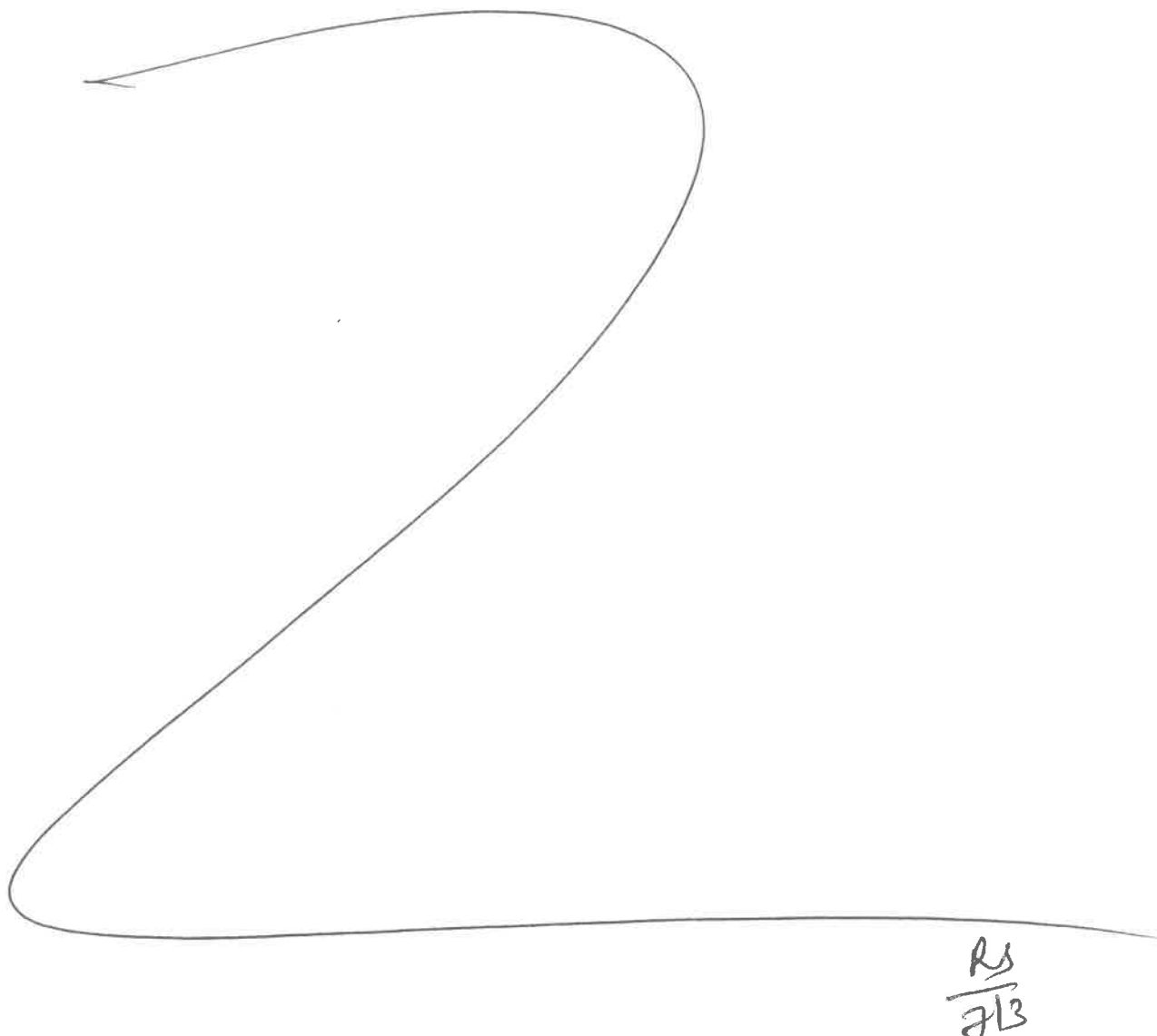
KD Bath Temperature: 60 °C **Envap Temperature:** 40 °C

Date / Time	Prepped Sample Relinquished By/Location	Received By/Location
7/3/25	RS (Ext Lab)	Y-Pelt+PUB
13235	Preparation Group	Analysis Group

Analytical Method: M3510C,3580A-Extraction Pesticide-17

Concentration Date: 07/03/2025

Sample ID	Client Sample ID	Test	g / mL	PH	Surr/Spike By:		Final Vol. (mL)	JarID	Comments	Prep Pos
					AddedBy	VerifiedBy				
PB168718BL	PBLK718	Pesticide-TCL	1000	6	RUPESH	ritesh	10			SEP-10
PB168718BS	PLCS718	Pesticide-TCL	1000	6	RUPESH	ritesh	10			11
PB168718BS D	PLCSD718	Pesticide-TCL	1000	6	RUPESH	ritesh	10			12
Q2458-10	FB-06272025	Pesticide-TCL	980	6	RUPESH	ritesh	10	I		13



8/14/25
8:52:38

WORKLIST(Hardcopy Internal Chain)

WorkList Name : Q2458P WorkList ID : 190540 Department : Extraction Date : 07-03-2025 08:51:40

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q2458-10	FB-06272025	Water	Herbicide	Cool 4 deg C	CAMP02	D51	06/27/2025	8151A
Q2458-10	FB-06272025	Water	Pesticide-TCL	Cool 4 deg C	CAMP02	D51	06/27/2025	8081B

Date/Time 7/13/25 8:52
Raw Sample Received by: RS (Ext 6-6)
Raw Sample Relinquished by: RS (Ext 6-6)

Date/Time 7/13/25 9:20
Raw Sample Received by: RS
Raw Sample Relinquished by: RS (Ext 6-6)



SHIPPING DOCUMENTS

CLIENT INFORMATION

REPORT TO BE SENT TO:

COMPANY: **CDM SMITH**
ADDRESS: **110 FIELDCREST AVE #8 6TH FLOOR**
CITY: **EDISON** STATE: **NJ** ZIP: **08837**
ATTENTION: **MARCIE ENCINAS**
PHONE: **7325904679** FAX:

CLIENT PROJECT INFORMATION

PROJECT NAME: **SOUTH RIVER WM REPLACEMENT**
PROJECT NO.: **302781** LOCATION: **SOUTH RIVER, NJ**
PROJECT MANAGER: **MARCIE ENCINAS**
e-mail: **ENCINASMA@CDMSMITH.COM**
PHONE: **7325904679** FAX:

CLIENT BILLING INFORMATION

BILL TO: **CDM SMITH** PO#: **7325904679**
ADDRESS: **110 FIELDCREST AVE #8 6TH FLOOR**
CITY: **EDISON** STATE: **NJ** ZIP: **08837**
ATTENTION: **MARCIE ENCINAS** PHONE: **7325904679**

ANALYSIS

DATA TURNAROUND INFORMATION

FAX (RUSH) _____ DAYS*
HARDCOPY (DATA PACKAGE): _____ DAYS*
EDD: _____ DAYS*
*TO BE APPROVED BY CHEMTECH
STANDARD HARDCOPY TURNAROUND TIME IS 10 BUSINESS

DATA DELIVERABLE INFORMATION

☐ Level 1 (Results Only) ☐ Level 4 (QC + Full Raw Data)
☒ Level 2 (Results + QC) ☐ NJ Reduced ☐ US EPA CLP
☐ Level 3 (Results + QC) ☐ NYS ASP A ☐ NYS ASP B
+ Raw Data ☐ Other _____
☐ EDD FORMAT _____

TCL VOC
TCL NVOC
PCB'S
TAL METALS
PESTICIDES
HERBICIDES
DRO/GRO

PRESERVATIVES

COMMENTS

ALLIANCE SAMPLE ID	PROJECT SAMPLE IDENTIFICATION	SAMPLE MATRIX	SAMPLE TYPE		SAMPLE COLLECTION		# OF BOTTLES										Specify Preservatives A-HCl D-NaOH B-HNO3 E-ICE C-H2SO4 F-OTHER
			COMP	GRAB	DATE	TIME		1	2	3	4	5	6	7	8	9	
1.	TP-70	S		X	6/26/25	1115	6	X	X	X	X	X	X	X			E
2.	TP-65	S		X	6/26/25	1200	6	X	X	X	X	X	X	X			E
3.	TP-68	S		X	6/27/25	0745	6	X	X	X	X	X	X	X			E
4.	TP-67	S		X	6/27/25	0805	6	X	X	X	X	X	X	X			E
5.	TP-66	S		X	6/27/25	0835	6	X	X	X	X	X	X	X			E
6.	TP-60	S		X	6/27/25	0910	6	X	X	X	X	X	X	X			E
7.	TP-62	S		X	6/27/25	1005	6	X	X	X	X	X	X	X			E
8.	TP-63	S		X	6/27/25	1055	6	X	X	X	X	X	X	X			E
9.	TP-54	S		X	6/27/25	1205	6	X	X	X	X	X	X	X			E
10.	FB-06272025	BAR		X	6/27/25	1300	9	X	X	X	X	X	X	X			E, A, B

SAMPLE CUSTODY MUST BE DOCUMENTED BELOW EACH TIME SAMPLES CHANGE POSSESSION INCLUDING COURIER DELIVERY

RELINQUISHED BY SAMPLER:	DATE/TIME:	RECEIVED BY:	Conditions of bottles or coolers at receipt: ☐ COMPLIANT ☐ NON COMPLIANT ☐ COOLER TEMP _____ °C
1. <i>[Signature]</i>	6/27/25/1600	<i>[Signature]</i> 1600	Comments: _____
RELINQUISHED BY SAMPLER:	DATE/TIME:	RECEIVED BY:	
2. <i>[Signature]</i>			
RELINQUISHED BY SAMPLER:	DATE/TIME:	RECEIVED BY:	Page _____ of _____
3. <i>[Signature]</i>	6-27-25		CLIENT: ☐ Hand Delivered ☐ Other
			Shipment Complete ☐ YES ☐ NO

From: Encinas, Marcie (Puskarik) <encinasma@cdmsmith.com>
Sent: Monday, June 30, 2025 12:44 PM
Subject: FW: sample id and sample not mentioned on COC

EXTERNAL EMAIL - This email was sent by a person from outside your organization. Exercise caution when clicking links, opening attachments or taking further action, before validating its authenticity.

Secured by Check Point

Thank you for reaching out - Please see below.

Would it be possible to also have excel results emailed with results compared to the current NJDEP SRS?

Marcie

From: Deepak Parmar <Deepak.Parmar@alliancetg.com>
Sent: Monday, June 30, 2025 10:24 AM
To: Encinas, Marcie (Puskarik) <encinasma@cdmsmith.com>
Cc: Mohammad Ahmed <mohammad.ahmed@alliancetg.com>; Yazmeen Gomez <Yazmeen.Gomez@alliancetg.com>
Subject: sample id and sample not mentioned on COC

You don't often get email from deepak.parmar@alliancetg.com. [Learn why this is important](#)

Good morning,

Samples received with 6/27/2025 shipment has below discrepancies

issue 1# sample (2) TP-65 mentioned on COC, however sample received as TP-55 let us know which sample ID to use ? **TP-55 is the correct label for the samples.**

Issue2# Lab receives two TB water samples but not mentioned on COC. Lab would like to know how to proceed with analysis ? **I believe those are temperature blanks, which do not require analysis. We are only analyzing trip blanks if the same also includes aqueous samples.**

Thanks & Regards,

Deepak Parmar
QA/QC
An Alliance Technical Group Company
Main: 908-789-8900

Address: 284 Sheffield St, Ste 1,
Mountainside, NJ 07092



Laboratory Certification

Certified By	License No.
CAS EPA CLP Contract	68HERH20D0011
Connecticut	PH-0830
DOD ELAP (ANAB)	L2219
Maine	2024021
Maryland	296
New Hampshire	255424 Rev 1
New Jersey	20012
New York	11376
Pennsylvania	68-00548
Soil Permit	525-24-234-08441
Texas	T104704488

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q2458	CAMP02	Order Date : 6/27/2025 4:22:00 PM	Project Mgr :
Client Name : CDM Smith		Project Name : South River WM Replacem	Report Type : Level 2 <i>all</i>
Client Contact : Marcie Ann Encinas		Receive DateTime : 6/27/2025 12:00:00 AM <i>10:55 PM</i> <i>all</i>	EDD Type : EXCEL NOCLEANUP
Invoice Name : CDM Smith		Purchase Order :	Hard Copy Date :
Invoice Contact : Marcie Ann Encinas			Date Signoff :

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
Q2458-01	TP-76	Solid	06/26/2025	11:15					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-02	TP- 65 <i>55 all</i>	Solid	06/26/2025	12:00					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-03	TP-68	Solid	06/27/2025	07: <i>45</i>					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-04	TP-67	Solid	06/27/2025	08:05					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-05	TP-66	Solid	06/27/2025	08:35					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-06	TP-60	Solid	06/27/2025	09:10					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-07	TP-62	Solid	06/27/2025	10:05					
					VOC-TCLVOA-10		8260D		10 Bus. Days
Q2458-08	TP-63	Solid	06/27/2025	10:55					

LOGIN REPORT/SAMPLE TRANSFER

Order ID : Q2458	CAMP02	Order Date : 6/27/2025 4:22:00 PM	Project Mgr :
Client Name : CDM Smith		Project Name : South River WM Replacem	Report Type : Level 2 <i>add</i>
Client Contact : Marcie Ann Encinas		Receive DateTime : 6/27/2025 12:00:00 AM <i>16:55 PM</i>	EDD Type : EXCEL NOCLEANUP
Invoice Name : CDM Smith		Purchase Order :	Hard Copy Date :
Invoice Contact : Marcie Ann Encinas			Date Signoff :

LAB ID	CLIENT ID	MATRIX	SAMPLE DATE	SAMPLE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
Q2458-09	TP-59	Solid	06/27/2025	12:05	VOC-TCLVOA-10		8260D	10 Bus. Days	
Q2458-10	FB-06272025	Water	06/27/2025	13:00	VOC-TCLVOA-10		8260D	10 Bus. Days	
					VOC-TCLVOA-10		8260-Low	10 Bus. Days	

Relinquished By : *[Signature]*

Date / Time : 6/30/25 0900

Received By : *[Signature]*

Date / Time : 6/30/25 9:00

Storage Area : VOA Refridgerator Room

SAMPLES RECEIVED ON 6/27/25 @ 1655
SAMPLES PLACED IN SM-REF-2

Ref #6
#2-2
Ref #4