



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

Hit Summary Sheet
SW-846

SDG No.: Q3495

Order ID: Q3495

Client: Remington & Vernick Engineers

Project ID: Maclearie Park

Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
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Client ID :

Total Concentration: 0.000

A
B
C
D
E
F
G
H
I
J
K
L



SAMPLE DATA



284 Sheffield Street, Mountainside, New Jersey 07092, Phone : 908 789 8900,
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Report of Analysis

Client:	Remington & Vernick Engineers	Date Collected:	10/27/25
Project:	Maclearie Park	Date Received:	10/29/25
Client Sample ID:	SB-1025-2	SDG No.:	Q3495
Lab Sample ID:	Q3495-02	Matrix:	SOIL
Analytical Method:	8081B	% Solid:	75.1
Sample Wt/Vol:	30.06 g	Final Vol:	10000 uL
Prep Method:	SW3541B	Test:	Pesticide-TCL
	Prep Date:	10/31/25	

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.17	U	1	0.17	2.30	ug/kg	10/31/25 13:33	PB170340
319-85-7	beta-BHC	0.24	U	1	0.24	2.30	ug/kg	10/31/25 13:33	PB170340
319-86-8	delta-BHC	0.52	U	1	0.52	2.30	ug/kg	10/31/25 13:33	PB170340
58-89-9	gamma-BHC (Lindane)	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
76-44-8	Heptachlor	0.16	U	1	0.16	2.30	ug/kg	10/31/25 13:33	PB170340
309-00-2	Aldrin	0.16	U	1	0.16	2.30	ug/kg	10/31/25 13:33	PB170340
1024-57-3	Heptachlor epoxide	0.25	U	1	0.25	2.30	ug/kg	10/31/25 13:33	PB170340
959-98-8	Endosulfan I	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
60-57-1	Dieldrin	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
72-55-9	4,4-DDE	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
72-20-8	Endrin	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
33213-65-9	Endosulfan II	0.39	U	1	0.39	2.30	ug/kg	10/31/25 13:33	PB170340
72-54-8	4,4-DDD	0.20	U	1	0.20	2.30	ug/kg	10/31/25 13:33	PB170340
1031-07-8	Endosulfan Sulfate	0.17	U	1	0.17	2.30	ug/kg	10/31/25 13:33	PB170340
50-29-3	4,4-DDT	0.19	U	1	0.19	2.30	ug/kg	10/31/25 13:33	PB170340
72-43-5	Methoxychlor	0.49	U	1	0.49	2.30	ug/kg	10/31/25 13:33	PB170340
53494-70-5	Endrin ketone	0.25	U	1	0.25	2.30	ug/kg	10/31/25 13:33	PB170340
7421-93-4	Endrin aldehyde	0.49	U	1	0.49	2.30	ug/kg	10/31/25 13:33	PB170340
5103-71-9	alpha-Chlordane	0.16	U	1	0.16	2.30	ug/kg	10/31/25 13:33	PB170340
5103-74-2	gamma-Chlordane	0.20	U	1	0.20	2.30	ug/kg	10/31/25 13:33	PB170340
8001-35-2	Toxaphene	7.20	U	1	7.20	43.9	ug/kg	10/31/25 13:33	PB170340
SURROGATES									
2051-24-3	Decachlorobiphenyl	16.1			20 - 144	81%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	19.7			19 - 148	99%	SPK: 20		

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



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Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/27/25
Project:	Maclearie Park		Date Received:	10/29/25
Client Sample ID:	TW-1025-1		SDG No.:	Q3495
Lab Sample ID:	Q3495-03		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	980 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0040	U	1	0.0040	0.051	ug/L	10/31/25 15:34	PB170348
319-85-7	beta-BHC	0.0050	U	1	0.0050	0.051	ug/L	10/31/25 15:34	PB170348
319-86-8	delta-BHC	0.011	U	1	0.011	0.051	ug/L	10/31/25 15:34	PB170348
58-89-9	gamma-BHC (Lindane)	0.0038	U	1	0.0038	0.051	ug/L	10/31/25 15:34	PB170348
76-44-8	Heptachlor	0.0028	U	1	0.0028	0.051	ug/L	10/31/25 15:34	PB170348
309-00-2	Aldrin	0.0037	U	1	0.0037	0.051	ug/L	10/31/25 15:34	PB170348
1024-57-3	Heptachlor epoxide	0.0098	U	1	0.0098	0.051	ug/L	10/31/25 15:34	PB170348
959-98-8	Endosulfan I	0.0032	U	1	0.0032	0.051	ug/L	10/31/25 15:34	PB170348
60-57-1	Dieldrin	0.0037	U	1	0.0037	0.051	ug/L	10/31/25 15:34	PB170348
72-55-9	4,4-DDE	0.0038	U	1	0.0038	0.051	ug/L	10/31/25 15:34	PB170348
72-20-8	Endrin	0.0033	U	1	0.0033	0.051	ug/L	10/31/25 15:34	PB170348
33213-65-9	Endosulfan II	0.0081	U	1	0.0081	0.051	ug/L	10/31/25 15:34	PB170348
72-54-8	4,4-DDD	0.0072	U	1	0.0072	0.051	ug/L	10/31/25 15:34	PB170348
1031-07-8	Endosulfan Sulfate	0.0038	U	1	0.0038	0.051	ug/L	10/31/25 15:34	PB170348
50-29-3	4,4-DDT	0.0036	U	1	0.0036	0.051	ug/L	10/31/25 15:34	PB170348
72-43-5	Methoxychlor	0.011	U	1	0.011	0.051	ug/L	10/31/25 15:34	PB170348
53494-70-5	Endrin ketone	0.0095	U	1	0.0095	0.051	ug/L	10/31/25 15:34	PB170348
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.051	ug/L	10/31/25 15:34	PB170348
5103-71-9	alpha-Chlordane	0.0036	U	1	0.0036	0.051	ug/L	10/31/25 15:34	PB170348
5103-74-2	gamma-Chlordane	0.0040	U	1	0.0040	0.051	ug/L	10/31/25 15:34	PB170348
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25 15:34	PB170348
SURROGATES									
2051-24-3	Decachlorobiphenyl	16.8			57 - 171	84%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	19.9			61 - 148	100%	SPK: 20		

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

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QC SUMMARY

Surrogate Summary

SDG No.: Q3495

Client: Remington & Vernick Engineers

Analytical Method: 8081B

Lab Sample ID	Client ID	Parameter	Column	Spike	Result	Recovery(%)	Qual	Limits(%)	
								Low	High
I.BLK-PD090647.D	PIBLK-PD090647.D	Decachlorobiphen	1	20	21.3	106		57	171
		Tetrachloro-m-xyl	1	20	19.4	97		61	148
		Decachlorobiphen	2	20	20.3	102		57	171
		Tetrachloro-m-xyl	2	20	20.0	100		61	148
I.BLK-PD090924.D	PIBLK-PD090924.D	Decachlorobiphen	1	20	22.3	111		57	171
		Tetrachloro-m-xyl	1	20	20.4	102		61	148
		Decachlorobiphen	2	20	28.7	143		57	171
		Tetrachloro-m-xyl	2	20	23.8	119		61	148
PB170348BL	PB170348BL	Decachlorobiphen	1	20	21.9	110		57	171
		Tetrachloro-m-xyl	1	20	18.8	94		61	148
		Decachlorobiphen	2	20	27.2	136		57	171
		Tetrachloro-m-xyl	2	20	22.4	112		61	148
Q3495-03	TW-1025-1	Decachlorobiphen	1	20	13.4	67		57	171
		Tetrachloro-m-xyl	1	20	16.6	83		61	148
		Decachlorobiphen	2	20	16.8	84		57	171
		Tetrachloro-m-xyl	2	20	19.9	100		61	148
I.BLK-PD090932.D	PIBLK-PD090932.D	Decachlorobiphen	1	20	21.9	110		57	171
		Tetrachloro-m-xyl	1	20	20.3	102		61	148
		Decachlorobiphen	2	20	28.1	140		57	171
		Tetrachloro-m-xyl	2	20	23.1	116		61	148
I.BLK-PD090948.D	PIBLK-PD090948.D	Decachlorobiphen	1	20	17.7	89		57	171
		Tetrachloro-m-xyl	1	20	19.2	96		61	148
		Decachlorobiphen	2	20	12.9	65		57	171
		Tetrachloro-m-xyl	2	20	19.9	99		61	148
PB170348BSD	PB170348BSD	Decachlorobiphen	1	20	25.0	125		57	171
		Tetrachloro-m-xyl	1	20	26.8	134		61	148
		Decachlorobiphen	2	20	21.3	106		57	171
		Tetrachloro-m-xyl	2	20	28.5	142		61	148
PB170348BS	PB170348BS	Decachlorobiphen	1	20	24.8	124		57	171
		Tetrachloro-m-xyl	1	20	27.6	138		61	148
		Decachlorobiphen	2	20	22.9	115		57	171
		Tetrachloro-m-xyl	2	20	28.8	144		61	148
I.BLK-PD090956.D	PIBLK-PD090956.D	Decachlorobiphen	1	20	19.2	96		57	171
		Tetrachloro-m-xyl	1	20	19.6	98		61	148
		Decachlorobiphen	2	20	17.8	89		57	171
		Tetrachloro-m-xyl	2	20	20.0	100		61	148
I.BLK-PL097636.D	PIBLK-PL097636.D	Decachlorobiphen	1	20	20.2	101		57	171
		Tetrachloro-m-xyl	1	20	18.0	90		61	148
		Decachlorobiphen	2	20	20.1	101		57	171
		Tetrachloro-m-xyl	2	20	18.7	94		61	148
I.BLK-PL097699.D	PIBLK-PL097699.D	Decachlorobiphen	1	20	22.7	113		57	171

Surrogate Summary

SDG No.: **Q3495**

Client: **Remington & Vernick Engineers**

Analytical Method: **8081B**

Lab Sample ID	Client ID	Parameter	Column	Spike	Result	Recovery(%)	Qual	Limits(%)	
								Low	High
I.BLK-PL097699.D	PIBLK-PL097699.D	Tetrachloro-m-xyl	1	20	23.6	118		61	148
		Decachlorobiphen	2	20	20.1	100		57	171
PB170340BL	PB170340BL	Tetrachloro-m-xyl	2	20	22.9	115		61	148
		Decachlorobiphen	1	20	18.9	94		20	144
		Tetrachloro-m-xyl	1	20	19.4	97		19	148
		Decachlorobiphen	2	20	15.9	80		20	144
		Tetrachloro-m-xyl	2	20	20.0	100		19	148
		Decachlorobiphen	1	20	23.4	117		20	144
PB170340BS	PB170340BS	Tetrachloro-m-xyl	1	20	23.7	118		19	148
		Decachlorobiphen	2	20	19.8	99		20	144
		Tetrachloro-m-xyl	2	20	23.1	116		19	148
		Decachlorobiphen	1	20	16.1	81		20	144
		Tetrachloro-m-xyl	1	20	19.7	99		19	148
		Decachlorobiphen	2	20	14.3	72		20	144
Q3495-02	SB-1025-2	Tetrachloro-m-xyl	2	20	18.9	94		19	148
		Decachlorobiphen	1	20	14.8	74		20	144
		Tetrachloro-m-xyl	1	20	16.8	84		19	148
		Decachlorobiphen	2	20	12.8	64		20	144
		Tetrachloro-m-xyl	2	20	16.7	83		19	148
		Decachlorobiphen	1	20	14.9	74		20	144
Q3495-02MS	SB-1025-2MS	Tetrachloro-m-xyl	1	20	17.0	85		19	148
		Decachlorobiphen	2	20	13.4	67		20	144
		Tetrachloro-m-xyl	2	20	17.2	86		19	148
		Decachlorobiphen	1	20	23.1	115		57	171
		Tetrachloro-m-xyl	1	20	23.0	115		61	148
		Decachlorobiphen	2	20	20.5	103		57	171
I.BLK-PL097714.D	PIBLK-PL097714.D	Tetrachloro-m-xyl	2	20	23.6	118		61	148

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q3495</u>	Analytical Method:	<u>8081B</u>
Client:	<u>Remington & Vernick Engineers</u>	DataFile :	<u>PL097707.D</u>

	Parameter	Spike	Sample		Units	Rec	Rec Qual	RPD	RPD Qual	Low	Limits		RPD
			Result								High		
Lab Sample ID:	Q3495-02MS (Column 1)	Client Sample ID:		SB-1025-2MS									
	alpha-BHC	22.16	0	22.1	ug/kg	100				60	144		
	beta-BHC	22.16	0	22.5	ug/kg	102				54	143		
	delta-BHC	22.16	0	23.0	ug/kg	104				29	151		
	gamma-BHC (Lindane)	22.16	0	22.1	ug/kg	100				61	140		
	Heptachlor	22.16	0	22.9	ug/kg	103				63	135		
	Aldrin	22.16	0	21.7	ug/kg	98				49	139		
	Heptachlor epoxide	22.16	0	21.1	ug/kg	95				41	156		
	Endosulfan I	22.16	0	21.4	ug/kg	97				56	142		
	Dieldrin	22.16	0	21.5	ug/kg	97				47	161		
	4,4'-DDE	22.16	0	22.6	ug/kg	102				55	136		
	Endrin	22.16	0	21.2	ug/kg	96				57	139		
	Endosulfan II	22.16	0	21.9	ug/kg	99				40	163		
	4,4'-DDD	22.16	0	23.3	ug/kg	105				47	163		
	Endosulfan sulfate	22.16	0	21.9	ug/kg	99				62	139		
	4,4'-DDT	22.16	0	22.1	ug/kg	100				51	146		
	Methoxychlor	22.16	0	21.5	ug/kg	97				54	136		
	Endrin ketone	22.16	0	21.7	ug/kg	98				60	129		
	Endrin aldehyde	22.16	0	21.6	ug/kg	97				59	132		
	alpha-Chlordane	22.16	0	21.4	ug/kg	97				39	166		
	gamma-Chlordane	22.16	0	21.9	ug/kg	99				44	175		
Lab Sample ID:	Q3495-02MS (Column 2)	Client Sample ID:		SB-1025-2MS									
	alpha-BHC	22.16	0	21.7	ug/kg	98				60	144		
	beta-BHC	22.16	0	20.7	ug/kg	93				54	143		
	delta-BHC	22.16	0	21.7	ug/kg	98				29	151		
	gamma-BHC (Lindane)	22.16	0	21.8	ug/kg	98				61	140		
	Heptachlor	22.16	0	21.4	ug/kg	97				63	135		
	Aldrin	22.16	0	21.8	ug/kg	98				49	139		
	Heptachlor epoxide	22.16	0	21.5	ug/kg	97				41	156		
	Endosulfan I	22.16	0	21.2	ug/kg	96				56	142		
	Dieldrin	22.16	0	21.6	ug/kg	97				47	161		
	4,4'-DDE	22.16	0	21.5	ug/kg	97				55	136		
	Endrin	22.16	0	22.1	ug/kg	100				57	139		
	Endosulfan II	22.16	0	21.0	ug/kg	95				40	163		
	4,4'-DDD	22.16	0	20.3	ug/kg	92				47	163		
	Endosulfan sulfate	22.16	0	20.2	ug/kg	91				62	139		

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q3495</u>	Analytical Method:	<u>8081B</u>
Client:	<u>Remington & Vernick Engineers</u>	DataFile :	<u>PL097707.D</u>

Parameter	Spike	Sample		Units	Rec	Rec Qual	RPD	RPD Qual	Low	Limits		RPD
		Result	Result							High		
4,4'-DDT	22.16	0	20.6	ug/kg	93				51	146		
Methoxychlor	22.16	0	19.6	ug/kg	88				54	136		
Endrin ketone	22.16	0	19.9	ug/kg	90				60	129		
Endrin aldehyde	22.16	0	19.2	ug/kg	87				59	132		
alpha-Chlordane	22.16	0	21.3	ug/kg	96				39	166		
gamma-Chlordane	22.16	0	21.1	ug/kg	95				44	175		

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q3495</u>	Analytical Method:	<u>8081B</u>
Client:	<u>Remington & Vernick Engineers</u>	DataFile :	<u>PL097708.D</u>

	Parameter	Spike	Sample		Units	Rec	Rec Qual	RPD	RPD Qual	Low	Limits		RPD
			Result								High		
Lab Sample ID:	Q3495-02MSD (Column 1)	Client Sample ID:		SB-1025-2MSD									
	alpha-BHC	22.18	0	22.7	ug/kg	102		2		60	144		20
	beta-BHC	22.18	0	23.0	ug/kg	104		2		54	143		20
	delta-BHC	22.18	0	23.6	ug/kg	106		2		29	151		20
	gamma-BHC (Lindane)	22.18	0	22.6	ug/kg	102		2		61	140		20
	Heptachlor	22.18	0	23.5	ug/kg	106		3		63	135		20
	Aldrin	22.18	0	22.2	ug/kg	100		2		49	139		20
	Heptachlor epoxide	22.18	0	21.6	ug/kg	97		2		41	156		20
	Endosulfan I	22.18	0	22.1	ug/kg	100		3		56	142		20
	Dieldrin	22.18	0	22.0	ug/kg	99		2		47	161		20
	4,4'-DDE	22.18	0	23.3	ug/kg	105		3		55	136		20
	Endrin	22.18	0	21.7	ug/kg	98		2		57	139		20
	Endosulfan II	22.18	0	22.4	ug/kg	101		2		40	163		20
	4,4'-DDD	22.18	0	23.8	ug/kg	107		2		47	163		20
	Endosulfan sulfate	22.18	0	22.3	ug/kg	101		2		62	139		20
	4,4'-DDT	22.18	0	22.4	ug/kg	101		1		51	146		20
	Methoxychlor	22.18	0	21.4	ug/kg	96		1		54	136		20
	Endrin ketone	22.18	0	22.3	ug/kg	101		3		60	129		20
	Endrin aldehyde	22.18	0	22.2	ug/kg	100		3		59	132		20
	alpha-Chlordane	22.18	0	22.1	ug/kg	100		3		39	166		20
	gamma-Chlordane	22.18	0	22.3	ug/kg	101		2		44	175		20
Lab Sample ID:	Q3495-02MSD (Column 2)	Client Sample ID:		SB-1025-2MSD									
	alpha-BHC	22.18	0	22.3	ug/kg	101		3		60	144		20
	beta-BHC	22.18	0	21.2	ug/kg	96		3		54	143		20
	delta-BHC	22.18	0	22.3	ug/kg	101		3		29	151		20
	gamma-BHC (Lindane)	22.18	0	22.4	ug/kg	101		3		61	140		20
	Heptachlor	22.18	0	21.7	ug/kg	98		1		63	135		20
	Aldrin	22.18	0	22.4	ug/kg	101		3		49	139		20
	Heptachlor epoxide	22.18	0	21.9	ug/kg	99		2		41	156		20
	Endosulfan I	22.18	0	21.7	ug/kg	98		2		56	142		20
	Dieldrin	22.18	0	22.0	ug/kg	99		2		47	161		20
	4,4'-DDE	22.18	0	22.0	ug/kg	99		2		55	136		20
	Endrin	22.18	0	22.5	ug/kg	101		1		57	139		20
	Endosulfan II	22.18	0	21.5	ug/kg	97		2		40	163		20
	4,4'-DDD	22.18	0	20.9	ug/kg	94		2		47	163		20
	Endosulfan sulfate	22.18	0	20.8	ug/kg	94		3		62	139		20

Matrix Spike/Matrix Spike Duplicate Summary

SW-846

SDG No.:	<u>Q3495</u>	Analytical Method:	<u>8081B</u>
Client:	<u>Remington & Vernick Engineers</u>	DataFile :	<u>PL097708.D</u>

Parameter	Spike	Sample		Units	Rec	Rec Qual	RPD	RPD Qual	Low	Limits	
		Result	Result							High	RPD
4,4'-DDT	22.18	0	21.0	ug/kg	95		2		51	146	20
Methoxychlor	22.18	0	20.0	ug/kg	90		2		54	136	20
Endrin ketone	22.18	0	20.5	ug/kg	92		2		60	129	20
Endrin aldehyde	22.18	0	19.7	ug/kg	89		2		59	132	20
alpha-Chlordane	22.18	0	21.7	ug/kg	98		2		39	166	20
gamma-Chlordane	22.18	0	21.5	ug/kg	97		2		44	175	20

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	Q3495	Analytical Method:	8081B
Client:	Remington & Vernick Engineers	Datafile :	PD090950.D

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD		Limits	
								Qual	Low	High	RPD
PB170348BSD (Column 1)	alpha-BHC	0.5	0.52	ug/L	104	0			85	130	20
	beta-BHC	0.5	0.49	ug/L	97	2			83	126	20
	delta-BHC	0.5	0.51	ug/L	103	0			69	141	20
	gamma-BHC (Lindane)	0.5	0.51	ug/L	103	0			82	129	20
	Heptachlor	0.5	0.63	ug/L	127	1			79	127	20
	Aldrin	0.5	0.51	ug/L	102	1			79	126	20
	Heptachlor epoxide	0.5	0.49	ug/L	97	4			81	124	20
	Endosulfan I	0.5	0.48	ug/L	97	1			85	122	20
	Dieldrin	0.5	0.49	ug/L	98	0			83	125	20
	4,4'-DDE	0.5	0.48	ug/L	97	1			80	127	20
	Endrin	0.5	0.47	ug/L	95	2			81	128	20
	Endosulfan II	0.5	0.48	ug/L	95	2			82	123	20
	4,4'-DDD	0.5	0.49	ug/L	98	1			77	131	20
	Endosulfan sulfate	0.5	0.45	ug/L	91	1			76	129	20
	4,4'-DDT	0.5	0.48	ug/L	95	1			80	133	20
	Methoxychlor	0.5	0.47	ug/L	94	2			78	108	20
	Endrin ketone	0.5	0.46	ug/L	91	1			80	131	20
	Endrin aldehyde	0.5	0.44	ug/L	89	1			82	127	20
	alpha-Chlordane	0.5	0.47	ug/L	93	4			82	125	20
	gamma-Chlordane	0.5	0.48	ug/L	97	3			82	125	20
	alpha-BHC	0.5	0.51	ug/L	102	1			85	130	20
	beta-BHC	0.5	0.50	ug/L	99	1			83	126	20
	delta-BHC	0.5	0.50	ug/L	101	1			69	141	20
	gamma-BHC (Lindane)	0.5	0.51	ug/L	101	1			82	129	20
PB170348BSD (Column 2)	Heptachlor	0.5	0.53	ug/L	106	1			79	127	20
	Aldrin	0.5	0.50	ug/L	101	0			79	126	20
	Heptachlor epoxide	0.5	0.50	ug/L	100	0			81	124	20
	Endosulfan I	0.5	0.49	ug/L	98	3			85	122	20
	Dieldrin	0.5	0.48	ug/L	96	3			83	125	20
	4,4'-DDE	0.5	0.48	ug/L	96	3			80	127	20
	Endrin	0.5	0.46	ug/L	92	4			81	128	20
	Endosulfan II	0.5	0.45	ug/L	91	3			82	123	20
	4,4'-DDD	0.5	0.47	ug/L	94	4			77	131	20
	Endosulfan sulfate	0.5	0.44	ug/L	88	4			76	129	20
	4,4'-DDT	0.5	0.46	ug/L	92	2			80	133	20
	Methoxychlor	0.5	0.43	ug/L	87	3			78	108	20
	Endrin ketone	0.5	0.41	ug/L	82	7			80	131	20
	Endrin aldehyde	0.5	0.45	ug/L	89	1			82	127	20
	alpha-Chlordane	0.5	0.49	ug/L	98	2			82	125	20
	gamma-Chlordane	0.5	0.49	ug/L	99	2			82	125	20

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	<u>Q3495</u>	Analytical Method:	<u>8081B</u>
Client:	<u>Remington & Vernick Engineers</u>	Datafile :	<u>PD090952.D</u>

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD	Limits		RPD
								Qual	Low	High	
PB170348BS (Column 1)	alpha-BHC	0.5	0.52	ug/L	104				85	130	
	beta-BHC	0.5	0.50	ug/L	99				83	126	
	delta-BHC	0.5	0.51	ug/L	103				69	141	
	gamma-BHC (Lindane)	0.5	0.51	ug/L	103				82	129	
	Heptachlor	0.5	0.63	ug/L	126				79	127	
	Aldrin	0.5	0.51	ug/L	103				79	126	
	Heptachlor epoxide	0.5	0.50	ug/L	101				81	124	
	Endosulfan I	0.5	0.49	ug/L	98				85	122	
	Dieldrin	0.5	0.49	ug/L	98				83	125	
	4,4'-DDE	0.5	0.48	ug/L	96				80	127	
	Endrin	0.5	0.48	ug/L	97				81	128	
	Endosulfan II	0.5	0.48	ug/L	97				82	123	
	4,4'-DDD	0.5	0.50	ug/L	99				77	131	
	Endosulfan sulfate	0.5	0.46	ug/L	92				76	129	
	4,4'-DDT	0.5	0.48	ug/L	96				80	133	
	Methoxychlor	0.5	0.48	ug/L	96				78	108	
	Endrin ketone	0.5	0.46	ug/L	92				80	131	
	Endrin aldehyde	0.5	0.45	ug/L	90				82	127	
	alpha-Chlordane	0.5	0.49	ug/L	97				82	125	
	gamma-Chlordane	0.5	0.50	ug/L	100				82	125	
PB170348BS (Column 2)	alpha-BHC	0.5	0.51	ug/L	103				85	130	
	beta-BHC	0.5	0.50	ug/L	100				83	126	
	delta-BHC	0.5	0.51	ug/L	102				69	141	
	gamma-BHC (Lindane)	0.5	0.51	ug/L	102				82	129	
	Heptachlor	0.5	0.54	ug/L	107				79	127	
	Aldrin	0.5	0.51	ug/L	101				79	126	
	Heptachlor epoxide	0.5	0.50	ug/L	100				81	124	
	Endosulfan I	0.5	0.50	ug/L	101				85	122	
	Dieldrin	0.5	0.49	ug/L	99				83	125	
	4,4'-DDE	0.5	0.50	ug/L	99				80	127	
	Endrin	0.5	0.48	ug/L	96				81	128	
	Endosulfan II	0.5	0.47	ug/L	94				82	123	
	4,4'-DDD	0.5	0.49	ug/L	98				77	131	
	Endosulfan sulfate	0.5	0.46	ug/L	92				76	129	
	4,4'-DDT	0.5	0.47	ug/L	94				80	133	
	Methoxychlor	0.5	0.45	ug/L	90				78	108	
	Endrin ketone	0.5	0.44	ug/L	88				80	131	
	Endrin aldehyde	0.5	0.45	ug/L	90				82	127	
	alpha-Chlordane	0.5	0.50	ug/L	100				82	125	
	gamma-Chlordane	0.5	0.50	ug/L	101				82	125	

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

SDG No.:	Q3495	Analytical Method:	8081B
Client:	Remington & Vernick Engineers	Datafile :	PL097705.D

Lab Sample ID	Parameter	Spike	Result	Units	Rec	RPD	Qual	RPD	Limits		RPD
								Qual	Low	High	
PB170340BS (Column 1)	alpha-BHC	16.65	17.0	ug/kg	102				84	123	
	beta-BHC	16.65	17.1	ug/kg	103				82	123	
	delta-BHC	16.65	17.4	ug/kg	105				83	126	
	gamma-BHC (Lindane)	16.65	17.0	ug/kg	102				83	125	
	Heptachlor	16.65	17.1	ug/kg	103				83	122	
	Aldrin	16.65	16.6	ug/kg	100				82	124	
	Heptachlor epoxide	16.65	16.0	ug/kg	96				83	120	
	Endosulfan I	16.65	16.1	ug/kg	97				81	124	
	Dieldrin	16.65	16.1	ug/kg	97				85	121	
	4,4'-DDE	16.65	16.8	ug/kg	101				81	123	
	Endrin	16.65	16.0	ug/kg	96				76	130	
	Endosulfan II	16.65	17.0	ug/kg	102				80	125	
	4,4'-DDD	16.65	18.0	ug/kg	108				80	131	
	Endosulfan sulfate	16.65	16.4	ug/kg	98				81	122	
	4,4'-DDT	16.65	16.3	ug/kg	98				70	129	
	Methoxychlor	16.65	15.2	ug/kg	91				60	119	
	Endrin ketone	16.65	16.8	ug/kg	101				77	132	
	Endrin aldehyde	16.65	16.3	ug/kg	98				79	124	
	alpha-Chlordane	16.65	16.2	ug/kg	97				84	120	
	gamma-Chlordane	16.65	16.5	ug/kg	99				83	122	
PB170340BS (Column 2)	alpha-BHC	16.65	16.9	ug/kg	102				84	123	
	beta-BHC	16.65	15.9	ug/kg	95				82	123	
	delta-BHC	16.65	16.7	ug/kg	100				83	126	
	gamma-BHC (Lindane)	16.65	16.9	ug/kg	102				83	125	
	Heptachlor	16.65	16.4	ug/kg	98				83	122	
	Aldrin	16.65	16.8	ug/kg	101				82	124	
	Heptachlor epoxide	16.65	16.3	ug/kg	98				83	120	
	Endosulfan I	16.65	15.8	ug/kg	95				81	124	
	Dieldrin	16.65	16.5	ug/kg	99				85	121	
	4,4'-DDE	16.65	16.3	ug/kg	98				81	123	
	Endrin	16.65	17.3	ug/kg	104				76	130	
	Endosulfan II	16.65	15.4	ug/kg	92				80	125	
	4,4'-DDD	16.65	16.1	ug/kg	97				80	131	
	Endosulfan sulfate	16.65	15.0	ug/kg	90				81	122	
	4,4'-DDT	16.65	14.9	ug/kg	89				70	129	
	Methoxychlor	16.65	14.2	ug/kg	85				60	119	
	Endrin ketone	16.65	16.1	ug/kg	97				77	132	
	Endrin aldehyde	16.65	14.2	ug/kg	85				79	124	
	alpha-Chlordane	16.65	16.1	ug/kg	97				84	120	
	gamma-Chlordane	16.65	16.0	ug/kg	96				83	122	

4C

PESTICIDE METHOD BLANK SUMMARY

Client ID

PB170340BL

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170340BL

Lab File ID: PL097704.D

Matrix: (soil/water) Solid

Extraction: (Type) SOXH

Sulfur Cleanup: (Y/N) N

Date Extracted: 10/31/2025

Date Analyzed (1): 10/31/2025

Date Analyzed (2): 10/31/2025

Time Analyzed (1): 13:06

Time Analyzed (2): 13:06

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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
PB170340BS	PB170340BS	PL097705.D	10/31/2025	10/31/2025
SB-1025-2	Q3495-02	PL097706.D	10/31/2025	10/31/2025
SB-1025-2MS	Q3495-02MS	PL097707.D	10/31/2025	10/31/2025
SB-1025-2MSD	Q3495-02MSD	PL097708.D	10/31/2025	10/31/2025

COMMENTS: _____

4C

PESTICIDE METHOD BLANK SUMMARY

Client ID

PB170348BL

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170348BL

Lab File ID: PD090927.D

Matrix: (soil/water) WATER

Extraction: (Type) SEPF

Sulfur Cleanup: (Y/N) N

Date Extracted: 10/31/2025

Date Analyzed (1): 10/31/2025

Date Analyzed (2): 10/31/2025

Time Analyzed (1): 14:37

Time Analyzed (2): 14:37

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
TW-1025-1	Q3495-03	PD090931.D	10/31/2025	10/31/2025
PB170348BSD	PB170348BSD	PD090950.D	11/04/2025	11/04/2025
PB170348BS	PB170348BS	PD090952.D	11/04/2025	11/04/2025

COMMENTS: _____



QC SAMPLE DATA



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

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	
Project:	Maclearie Park		Date Received:	
Client Sample ID:	PB170340BL		SDG No.:	Q3495
Lab Sample ID:	PB170340BL		Matrix:	SOIL
Analytical Method:	8081B		% Solid:	100
Sample Wt/Vol:	30.02 g	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	SW3541B	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.13	U	1	0.13	1.70	ug/kg	10/31/25 13:06	PB170340
319-85-7	beta-BHC	0.18	U	1	0.18	1.70	ug/kg	10/31/25 13:06	PB170340
319-86-8	delta-BHC	0.39	U	1	0.39	1.70	ug/kg	10/31/25 13:06	PB170340
58-89-9	gamma-BHC (Lindane)	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
76-44-8	Heptachlor	0.12	U	1	0.12	1.70	ug/kg	10/31/25 13:06	PB170340
309-00-2	Aldrin	0.12	U	1	0.12	1.70	ug/kg	10/31/25 13:06	PB170340
1024-57-3	Heptachlor epoxide	0.19	U	1	0.19	1.70	ug/kg	10/31/25 13:06	PB170340
959-98-8	Endosulfan I	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
60-57-1	Dieldrin	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
72-55-9	4,4-DDE	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
72-20-8	Endrin	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
33213-65-9	Endosulfan II	0.29	U	1	0.29	1.70	ug/kg	10/31/25 13:06	PB170340
72-54-8	4,4-DDD	0.15	U	1	0.15	1.70	ug/kg	10/31/25 13:06	PB170340
1031-07-8	Endosulfan Sulfate	0.13	U	1	0.13	1.70	ug/kg	10/31/25 13:06	PB170340
50-29-3	4,4-DDT	0.14	U	1	0.14	1.70	ug/kg	10/31/25 13:06	PB170340
72-43-5	Methoxychlor	0.37	U	1	0.37	1.70	ug/kg	10/31/25 13:06	PB170340
53494-70-5	Endrin ketone	0.19	U	1	0.19	1.70	ug/kg	10/31/25 13:06	PB170340
7421-93-4	Endrin aldehyde	0.37	U	1	0.37	1.70	ug/kg	10/31/25 13:06	PB170340
5103-71-9	alpha-Chlordane	0.12	U	1	0.12	1.70	ug/kg	10/31/25 13:06	PB170340
5103-74-2	gamma-Chlordane	0.15	U	1	0.15	1.70	ug/kg	10/31/25 13:06	PB170340
8001-35-2	Toxaphene	5.40	U	1	5.40	33.0	ug/kg	10/31/25 13:06	PB170340
SURROGATES									
2051-24-3	Decachlorobiphenyl	18.9			20 - 144	94%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	20.0			19 - 148	100%	SPK: 20		

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



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

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	
Project:	Maclearie Park		Date Received:	
Client Sample ID:	PB170348BL		SDG No.:	Q3495
Lab Sample ID:	PB170348BL		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/31/25 14:37	PB170348
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/31/25 14:37	PB170348
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/31/25 14:37	PB170348
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/31/25 14:37	PB170348
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/31/25 14:37	PB170348
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25 14:37	PB170348
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/31/25 14:37	PB170348
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/31/25 14:37	PB170348
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25 14:37	PB170348
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/31/25 14:37	PB170348
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/31/25 14:37	PB170348
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/31/25 14:37	PB170348
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/31/25 14:37	PB170348
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/31/25 14:37	PB170348
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/31/25 14:37	PB170348
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/31/25 14:37	PB170348
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/31/25 14:37	PB170348
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/31/25 14:37	PB170348
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/31/25 14:37	PB170348
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/31/25 14:37	PB170348
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25 14:37	PB170348
SURROGATES									
2051-24-3	Decachlorobiphenyl	27.2			57 - 171	136%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	22.4			61 - 148	112%	SPK: 20		

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

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/17/25
Project:	Maclearie Park		Date Received:	10/17/25
Client Sample ID:	PIBLK-PD090647.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PD090647.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/17/25	pd101725
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/17/25	pd101725
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/17/25	pd101725
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/17/25	pd101725
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/17/25	pd101725
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/17/25	pd101725
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/17/25	pd101725
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/17/25	pd101725
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/17/25	pd101725
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/17/25	pd101725
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/17/25	pd101725
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/17/25	pd101725
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/17/25	pd101725
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/17/25	pd101725
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/17/25	pd101725
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/17/25	pd101725
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/17/25	pd101725
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/17/25	pd101725
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/17/25	pd101725
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/17/25	pd101725
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/17/25	pd101725
SURROGATES									
2051-24-3	Decachlorobiphenyl	21.3			57 - 171	106%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	20.0			61 - 148	100%	SPK: 20		

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

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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

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/31/25
Project:	Maclearie Park		Date Received:	10/31/25
Client Sample ID:	PIBLK-PD090924.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PD090924.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pd103125
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/31/25	pd103125
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/31/25	pd103125
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pd103125
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/31/25	pd103125
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/31/25	pd103125
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pd103125
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/31/25	pd103125
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/31/25	pd103125
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/31/25	pd103125
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pd103125
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/31/25	pd103125
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pd103125
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pd103125
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25	pd103125
SURROGATES									
2051-24-3	Decachlorobiphenyl	28.7			57 - 171	143%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	23.8			61 - 148	119%	SPK: 20		

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Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/31/25
Project:	Maclearie Park		Date Received:	10/31/25
Client Sample ID:	PIBLK-PD090932.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PD090932.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pd103125
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/31/25	pd103125
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/31/25	pd103125
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pd103125
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/31/25	pd103125
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/31/25	pd103125
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pd103125
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/31/25	pd103125
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/31/25	pd103125
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/31/25	pd103125
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pd103125
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pd103125
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/31/25	pd103125
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/31/25	pd103125
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pd103125
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pd103125
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25	pd103125
SURROGATES									
2051-24-3	Decachlorobiphenyl	28.1			57 - 171	140%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	23.1			61 - 148	116%	SPK: 20		

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() = Laboratory InHouse Limit

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	11/04/25
Project:	Maclearie Park		Date Received:	11/04/25
Client Sample ID:	PIBLK-PD090948.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PD090948.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	11/04/25	pd110425
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	11/04/25	pd110425
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	11/04/25	pd110425
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	11/04/25	pd110425
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	11/04/25	pd110425
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	11/04/25	pd110425
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	11/04/25	pd110425
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	11/04/25	pd110425
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	11/04/25	pd110425
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	11/04/25	pd110425
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	11/04/25	pd110425
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	11/04/25	pd110425
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	11/04/25	pd110425
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	11/04/25	pd110425
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	11/04/25	pd110425
SURROGATES									
2051-24-3	Decachlorobiphenyl	17.7			57 - 171	89%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	19.9			61 - 148	99%	SPK: 20		

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Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	11/04/25
Project:	Maclearie Park		Date Received:	11/04/25
Client Sample ID:	PIBLK-PD090956.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PD090956.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	11/04/25	pd110425
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	11/04/25	pd110425
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	11/04/25	pd110425
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	11/04/25	pd110425
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	11/04/25	pd110425
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	11/04/25	pd110425
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	11/04/25	pd110425
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	11/04/25	pd110425
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	11/04/25	pd110425
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	11/04/25	pd110425
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	11/04/25	pd110425
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	11/04/25	pd110425
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	11/04/25	pd110425
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	11/04/25	pd110425
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	11/04/25	pd110425
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	11/04/25	pd110425
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	11/04/25	pd110425
SURROGATES									
2051-24-3	Decachlorobiphenyl	19.2			57 - 171	96%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	20.0			61 - 148	100%	SPK: 20		

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Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/28/25
Project:	Maclearie Park		Date Received:	10/28/25
Client Sample ID:	PIBLK-PL097636.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PL097636.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/28/25	PL102825
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/28/25	PL102825
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/28/25	PL102825
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/28/25	PL102825
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/28/25	PL102825
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/28/25	PL102825
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/28/25	PL102825
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/28/25	PL102825
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/28/25	PL102825
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/28/25	PL102825
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/28/25	PL102825
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/28/25	PL102825
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/28/25	PL102825
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/28/25	PL102825
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/28/25	PL102825
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/28/25	PL102825
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/28/25	PL102825
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/28/25	PL102825
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/28/25	PL102825
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/28/25	PL102825
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/28/25	PL102825
SURROGATES									
2051-24-3	Decachlorobiphenyl	20.2			57 - 171	101%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	18.7			61 - 148	94%	SPK: 20		

U = Not Detected

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Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/31/25
Project:	Maclearie Park		Date Received:	10/31/25
Client Sample ID:	PIBLK-PL097699.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PL097699.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pl103125
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/31/25	pl103125
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/31/25	pl103125
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pl103125
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/31/25	pl103125
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/31/25	pl103125
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pl103125
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/31/25	pl103125
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/31/25	pl103125
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/31/25	pl103125
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pl103125
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/31/25	pl103125
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pl103125
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pl103125
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25	pl103125
SURROGATES									
2051-24-3	Decachlorobiphenyl	22.7			57 - 171	113%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	23.6			61 - 148	118%	SPK: 20		

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

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J = Estimated Value

B = Analyte Found in Associated Method Blank

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

() = Laboratory InHouse Limit

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	10/31/25
Project:	Maclearie Park		Date Received:	10/31/25
Client Sample ID:	PIBLK-PL097714.D		SDG No.:	Q3495
Lab Sample ID:	I.BLK-PL097714.D		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date:		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pl103125
319-85-7	beta-BHC	0.0049	U	1	0.0049	0.050	ug/L	10/31/25	pl103125
319-86-8	delta-BHC	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
58-89-9	gamma-BHC (Lindane)	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
76-44-8	Heptachlor	0.0027	U	1	0.0027	0.050	ug/L	10/31/25	pl103125
309-00-2	Aldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pl103125
1024-57-3	Heptachlor epoxide	0.0096	U	1	0.0096	0.050	ug/L	10/31/25	pl103125
959-98-8	Endosulfan I	0.0031	U	1	0.0031	0.050	ug/L	10/31/25	pl103125
60-57-1	Dieldrin	0.0036	U	1	0.0036	0.050	ug/L	10/31/25	pl103125
72-55-9	4,4-DDE	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
72-20-8	Endrin	0.0032	U	1	0.0032	0.050	ug/L	10/31/25	pl103125
33213-65-9	Endosulfan II	0.0079	U	1	0.0079	0.050	ug/L	10/31/25	pl103125
72-54-8	4,4-DDD	0.0071	U	1	0.0071	0.050	ug/L	10/31/25	pl103125
1031-07-8	Endosulfan Sulfate	0.0037	U	1	0.0037	0.050	ug/L	10/31/25	pl103125
50-29-3	4,4-DDT	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pl103125
72-43-5	Methoxychlor	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
53494-70-5	Endrin ketone	0.0093	U	1	0.0093	0.050	ug/L	10/31/25	pl103125
7421-93-4	Endrin aldehyde	0.011	U	1	0.011	0.050	ug/L	10/31/25	pl103125
5103-71-9	alpha-Chlordane	0.0035	U	1	0.0035	0.050	ug/L	10/31/25	pl103125
5103-74-2	gamma-Chlordane	0.0039	U	1	0.0039	0.050	ug/L	10/31/25	pl103125
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	10/31/25	pl103125
SURROGATES									
2051-24-3	Decachlorobiphenyl	23.1			57 - 171	115%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	23.6			61 - 148	118%	SPK: 20		

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() = Laboratory InHouse Limit



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Fax : 908 789 8922

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	
Project:	Maclearie Park		Date Received:	
Client Sample ID:	PB170340BS		SDG No.:	Q3495
Lab Sample ID:	PB170340BS		Matrix:	SOIL
Analytical Method:	8081B		% Solid:	100
Sample Wt/Vol:	30.03 g	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	SW3541B	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	17.0		1	0.13	1.70	ug/kg	10/31/25 13:19	PB170340
319-85-7	beta-BHC	17.1		1	0.18	1.70	ug/kg	10/31/25 13:19	PB170340
319-86-8	delta-BHC	17.4		1	0.39	1.70	ug/kg	10/31/25 13:19	PB170340
58-89-9	gamma-BHC (Lindane)	17.0		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
76-44-8	Heptachlor	17.1		1	0.12	1.70	ug/kg	10/31/25 13:19	PB170340
309-00-2	Aldrin	16.8		1	0.12	1.70	ug/kg	10/31/25 13:19	PB170340
1024-57-3	Heptachlor epoxide	16.3		1	0.19	1.70	ug/kg	10/31/25 13:19	PB170340
959-98-8	Endosulfan I	16.1		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
60-57-1	Dieldrin	16.5		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
72-55-9	4,4-DDE	16.8		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
72-20-8	Endrin	17.3		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
33213-65-9	Endosulfan II	17.0		1	0.29	1.70	ug/kg	10/31/25 13:19	PB170340
72-54-8	4,4-DDD	18.0		1	0.15	1.70	ug/kg	10/31/25 13:19	PB170340
1031-07-8	Endosulfan Sulfate	16.4		1	0.13	1.70	ug/kg	10/31/25 13:19	PB170340
50-29-3	4,4-DDT	16.3		1	0.14	1.70	ug/kg	10/31/25 13:19	PB170340
72-43-5	Methoxychlor	15.2		1	0.37	1.70	ug/kg	10/31/25 13:19	PB170340
53494-70-5	Endrin ketone	16.8		1	0.19	1.70	ug/kg	10/31/25 13:19	PB170340
7421-93-4	Endrin aldehyde	16.3		1	0.37	1.70	ug/kg	10/31/25 13:19	PB170340
5103-71-9	alpha-Chlordane	16.2		1	0.12	1.70	ug/kg	10/31/25 13:19	PB170340
5103-74-2	gamma-Chlordane	16.5		1	0.15	1.70	ug/kg	10/31/25 13:19	PB170340
8001-35-2	Toxaphene	5.40	U	1	5.40	33.0	ug/kg	10/31/25 13:19	PB170340
SURROGATES									
2051-24-3	Decachlorobiphenyl	23.4			20 - 144	117%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	23.7			19 - 148	118%	SPK: 20		

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

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	
Project:	Maclearie Park		Date Received:	
Client Sample ID:	PB170348BS		SDG No.:	Q3495
Lab Sample ID:	PB170348BS		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.52		1	0.0039	0.050	ug/L	11/04/25 14:22	PB170348
319-85-7	beta-BHC	0.50		1	0.0049	0.050	ug/L	11/04/25 14:22	PB170348
319-86-8	delta-BHC	0.51		1	0.011	0.050	ug/L	11/04/25 14:22	PB170348
58-89-9	gamma-BHC (Lindane)	0.51		1	0.0037	0.050	ug/L	11/04/25 14:22	PB170348
76-44-8	Heptachlor	0.63		1	0.0027	0.050	ug/L	11/04/25 14:22	PB170348
309-00-2	Aldrin	0.51		1	0.0036	0.050	ug/L	11/04/25 14:22	PB170348
1024-57-3	Heptachlor epoxide	0.50		1	0.0096	0.050	ug/L	11/04/25 14:22	PB170348
959-98-8	Endosulfan I	0.50		1	0.0031	0.050	ug/L	11/04/25 14:22	PB170348
60-57-1	Dieldrin	0.49		1	0.0036	0.050	ug/L	11/04/25 14:22	PB170348
72-55-9	4,4-DDE	0.50		1	0.0037	0.050	ug/L	11/04/25 14:22	PB170348
72-20-8	Endrin	0.48		1	0.0032	0.050	ug/L	11/04/25 14:22	PB170348
33213-65-9	Endosulfan II	0.48		1	0.0079	0.050	ug/L	11/04/25 14:22	PB170348
72-54-8	4,4-DDD	0.50		1	0.0071	0.050	ug/L	11/04/25 14:22	PB170348
1031-07-8	Endosulfan Sulfate	0.46		1	0.0037	0.050	ug/L	11/04/25 14:22	PB170348
50-29-3	4,4-DDT	0.48		1	0.0035	0.050	ug/L	11/04/25 14:22	PB170348
72-43-5	Methoxychlor	0.48		1	0.011	0.050	ug/L	11/04/25 14:22	PB170348
53494-70-5	Endrin ketone	0.46		1	0.0093	0.050	ug/L	11/04/25 14:22	PB170348
7421-93-4	Endrin aldehyde	0.45		1	0.011	0.050	ug/L	11/04/25 14:22	PB170348
5103-71-9	alpha-Chlordane	0.50		1	0.0035	0.050	ug/L	11/04/25 14:22	PB170348
5103-74-2	gamma-Chlordane	0.50		1	0.0039	0.050	ug/L	11/04/25 14:22	PB170348
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	11/04/25 14:22	PB170348
SURROGATES									
2051-24-3	Decachlorobiphenyl	24.8			57 - 171	124%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	28.8			61 - 148	144%	SPK: 20		

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Fax : 908 789 8922

Report of Analysis

Client:	Remington & Vernick Engineers		Date Collected:	
Project:	Maclearie Park		Date Received:	
Client Sample ID:	PB170348BSD		SDG No.:	Q3495
Lab Sample ID:	PB170348BSD		Matrix:	WATER
Analytical Method:	8081B		% Solid:	0
Sample Wt/Vol:	1000 mL	Final Vol: 10000 uL	Test:	Pesticide-TCL
Prep Method:	3510C	Prep Date: 10/31/25		

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	0.52		1	0.0039	0.050	ug/L	11/04/25 13:52	PB170348
319-85-7	beta-BHC	0.50		1	0.0049	0.050	ug/L	11/04/25 13:52	PB170348
319-86-8	delta-BHC	0.51		1	0.011	0.050	ug/L	11/04/25 13:52	PB170348
58-89-9	gamma-BHC (Lindane)	0.51		1	0.0037	0.050	ug/L	11/04/25 13:52	PB170348
76-44-8	Heptachlor	0.63		1	0.0027	0.050	ug/L	11/04/25 13:52	PB170348
309-00-2	Aldrin	0.51		1	0.0036	0.050	ug/L	11/04/25 13:52	PB170348
1024-57-3	Heptachlor epoxide	0.50		1	0.0096	0.050	ug/L	11/04/25 13:52	PB170348
959-98-8	Endosulfan I	0.49		1	0.0031	0.050	ug/L	11/04/25 13:52	PB170348
60-57-1	Dieldrin	0.49		1	0.0036	0.050	ug/L	11/04/25 13:52	PB170348
72-55-9	4,4-DDE	0.48		1	0.0037	0.050	ug/L	11/04/25 13:52	PB170348
72-20-8	Endrin	0.47		1	0.0032	0.050	ug/L	11/04/25 13:52	PB170348
33213-65-9	Endosulfan II	0.48		1	0.0079	0.050	ug/L	11/04/25 13:52	PB170348
72-54-8	4,4-DDD	0.49		1	0.0071	0.050	ug/L	11/04/25 13:52	PB170348
1031-07-8	Endosulfan Sulfate	0.45		1	0.0037	0.050	ug/L	11/04/25 13:52	PB170348
50-29-3	4,4-DDT	0.48		1	0.0035	0.050	ug/L	11/04/25 13:52	PB170348
72-43-5	Methoxychlor	0.47		1	0.011	0.050	ug/L	11/04/25 13:52	PB170348
53494-70-5	Endrin ketone	0.46		1	0.0093	0.050	ug/L	11/04/25 13:52	PB170348
7421-93-4	Endrin aldehyde	0.45		1	0.011	0.050	ug/L	11/04/25 13:52	PB170348
5103-71-9	alpha-Chlordane	0.49		1	0.0035	0.050	ug/L	11/04/25 13:52	PB170348
5103-74-2	gamma-Chlordane	0.49		1	0.0039	0.050	ug/L	11/04/25 13:52	PB170348
8001-35-2	Toxaphene	0.17	U	1	0.17	1.00	ug/L	11/04/25 13:52	PB170348
SURROGATES									
2051-24-3	Decachlorobiphenyl	25.0			57 - 171	125%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	28.5			61 - 148	142%	SPK: 20		

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Report of Analysis

Client: Remington & Vernick Engineers	Date Collected: 10/27/25	
Project: Maclearie Park	Date Received: 10/29/25	
Client Sample ID: SB-1025-2MS	SDG No.: Q3495	
Lab Sample ID: Q3495-02MS	Matrix: SOIL	
Analytical Method: 8081B	% Solid: 75.1	
Sample Wt/Vol: 30.04 g	Test: Pesticide-TCL	Final Vol: 10000 uL
Prep Method: SW3541B		Prep Date: 10/31/25

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	22.1		1	0.17	2.30	ug/kg	10/31/25 13:47	PB170340
319-85-7	beta-BHC	22.5		1	0.24	2.30	ug/kg	10/31/25 13:47	PB170340
319-86-8	delta-BHC	23.0		1	0.52	2.30	ug/kg	10/31/25 13:47	PB170340
58-89-9	gamma-BHC (Lindane)	22.1		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
76-44-8	Heptachlor	22.9		1	0.16	2.30	ug/kg	10/31/25 13:47	PB170340
309-00-2	Aldrin	21.8		1	0.16	2.30	ug/kg	10/31/25 13:47	PB170340
1024-57-3	Heptachlor epoxide	21.5		1	0.25	2.30	ug/kg	10/31/25 13:47	PB170340
959-98-8	Endosulfan I	21.4		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
60-57-1	Dieldrin	21.6		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
72-55-9	4,4-DDE	22.6		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
72-20-8	Endrin	22.1		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
33213-65-9	Endosulfan II	21.9		1	0.39	2.30	ug/kg	10/31/25 13:47	PB170340
72-54-8	4,4-DDD	23.3		1	0.20	2.30	ug/kg	10/31/25 13:47	PB170340
1031-07-8	Endosulfan Sulfate	21.9		1	0.17	2.30	ug/kg	10/31/25 13:47	PB170340
50-29-3	4,4-DDT	22.1		1	0.19	2.30	ug/kg	10/31/25 13:47	PB170340
72-43-5	Methoxychlor	21.5		1	0.49	2.30	ug/kg	10/31/25 13:47	PB170340
53494-70-5	Endrin ketone	21.7		1	0.25	2.30	ug/kg	10/31/25 13:47	PB170340
7421-93-4	Endrin aldehyde	21.6		1	0.49	2.30	ug/kg	10/31/25 13:47	PB170340
5103-71-9	alpha-Chlordane	21.4		1	0.16	2.30	ug/kg	10/31/25 13:47	PB170340
5103-74-2	gamma-Chlordane	21.9		1	0.20	2.30	ug/kg	10/31/25 13:47	PB170340
8001-35-2	Toxaphene	7.20	U	1	7.20	43.9	ug/kg	10/31/25 13:47	PB170340
SURROGATES									
2051-24-3	Decachlorobiphenyl	14.8			20 - 144	74%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	16.8			19 - 148	84%	SPK: 20		

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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

Report of Analysis

Client: Remington & Vernick Engineers	Date Collected: 10/27/25	
Project: Maclearie Park	Date Received: 10/29/25	
Client Sample ID: SB-1025-2MSD	SDG No.: Q3495	
Lab Sample ID: Q3495-02MSD	Matrix: SOIL	
Analytical Method: 8081B	% Solid: 75.1	
Sample Wt/Vol: 30.02 g	Test: Pesticide-TCL	Final Vol: 10000 uL
Prep Method: SW3541B		Prep Date: 10/31/25

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
319-84-6	alpha-BHC	22.7		1	0.17	2.30	ug/kg	10/31/25 14:00	PB170340
319-85-7	beta-BHC	23.0		1	0.24	2.30	ug/kg	10/31/25 14:00	PB170340
319-86-8	delta-BHC	23.6		1	0.52	2.30	ug/kg	10/31/25 14:00	PB170340
58-89-9	gamma-BHC (Lindane)	22.6		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
76-44-8	Heptachlor	23.5		1	0.16	2.30	ug/kg	10/31/25 14:00	PB170340
309-00-2	Aldrin	22.4		1	0.16	2.30	ug/kg	10/31/25 14:00	PB170340
1024-57-3	Heptachlor epoxide	21.9		1	0.25	2.30	ug/kg	10/31/25 14:00	PB170340
959-98-8	Endosulfan I	22.1		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
60-57-1	Dieldrin	22.0		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
72-55-9	4,4-DDE	23.3		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
72-20-8	Endrin	22.5		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
33213-65-9	Endosulfan II	22.4		1	0.39	2.30	ug/kg	10/31/25 14:00	PB170340
72-54-8	4,4-DDD	23.8		1	0.20	2.30	ug/kg	10/31/25 14:00	PB170340
1031-07-8	Endosulfan Sulfate	22.3		1	0.17	2.30	ug/kg	10/31/25 14:00	PB170340
50-29-3	4,4-DDT	22.4		1	0.19	2.30	ug/kg	10/31/25 14:00	PB170340
72-43-5	Methoxychlor	21.4		1	0.49	2.30	ug/kg	10/31/25 14:00	PB170340
53494-70-5	Endrin ketone	22.3		1	0.25	2.30	ug/kg	10/31/25 14:00	PB170340
7421-93-4	Endrin aldehyde	22.2		1	0.49	2.30	ug/kg	10/31/25 14:00	PB170340
5103-71-9	alpha-Chlordane	22.1		1	0.16	2.30	ug/kg	10/31/25 14:00	PB170340
5103-74-2	gamma-Chlordane	22.3		1	0.20	2.30	ug/kg	10/31/25 14:00	PB170340
8001-35-2	Toxaphene	7.20	U	1	7.20	43.9	ug/kg	10/31/25 14:00	PB170340
SURROGATES									
2051-24-3	Decachlorobiphenyl	14.9			20 - 144	74%	SPK: 20		
877-09-8	Tetrachloro-m-xylene	17.2			19 - 148	86%	SPK: 20		

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



CALIBRATION SUMMARY

RETENTION TIMES OF INITIAL CALIBRATION

Lab Name:	<u>Alliance</u>	Contract:	<u>REMI02</u>
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q3495</u>
Instrument ID:	<u>ECD_D</u>	Calibration Date(s):	<u>10/17/2025</u> <u>10/17/2025</u>
		Calibration Times:	<u>11:34</u> <u>12:28</u>

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

LAB FILE ID:	RT 100 = <u>PD090650.D</u>	RT 075 = <u>PD090651.D</u>
RT 050 = <u>PD090652.D</u>	RT 025 = <u>PD090653.D</u>	RT 005 = <u>PD090654.D</u>

COMPOUND	RT 100	RT 075	RT 050	RT 025	RT 005	MEAN RT	RT WINDOW	
							FROM	TO
4,4'-DDD	6.70	6.70	6.70	6.70	6.70	6.70	6.60	6.80
4,4'-DDE	6.19	6.19	6.19	6.19	6.19	6.19	6.09	6.29
4,4'-DDT	7.02	7.02	7.02	7.02	7.02	7.02	6.92	7.12
Aldrin	5.26	5.27	5.26	5.26	5.26	5.26	5.16	5.36
alpha-BHC	3.99	3.99	3.99	3.99	3.99	3.99	3.89	4.09
alpha-Chlordane	6.02	6.02	6.02	6.02	6.02	6.02	5.92	6.12
beta-BHC	4.51	4.51	4.51	4.51	4.51	4.51	4.41	4.61
Decachlorobiphenyl	9.07	9.07	9.06	9.07	9.07	9.07	8.97	9.17
delta-BHC	4.76	4.76	4.76	4.76	4.76	4.76	4.66	4.86
Dieldrin	6.34	6.34	6.34	6.34	6.34	6.34	6.24	6.44
Endosulfan I	6.07	6.07	6.07	6.07	6.07	6.07	5.97	6.17
Endosulfan II	6.78	6.78	6.78	6.78	6.78	6.78	6.68	6.88
Endosulfan sulfate	7.15	7.15	7.15	7.14	7.14	7.14	7.04	7.24
Endrin	6.57	6.57	6.57	6.57	6.57	6.57	6.47	6.67
Endrin aldehyde	6.91	6.91	6.91	6.91	6.91	6.91	6.81	7.01
Endrin ketone	7.63	7.63	7.63	7.63	7.62	7.62	7.52	7.72
gamma-BHC (Lindane)	4.32	4.33	4.33	4.32	4.32	4.32	4.22	4.42
gamma-Chlordane	5.94	5.94	5.94	5.94	5.94	5.94	5.84	6.04
Heptachlor	4.92	4.92	4.92	4.92	4.92	4.92	4.82	5.02
Heptachlor epoxide	5.69	5.69	5.68	5.68	5.68	5.68	5.58	5.78
Methoxychlor	7.49	7.49	7.49	7.49	7.49	7.49	7.39	7.59
Tetrachloro-m-xylene	3.54	3.55	3.55	3.54	3.54	3.54	3.44	3.64

RETENTION TIMES OF INITIAL CALIBRATION

Lab Name:	<u>Alliance</u>	Contract:	<u>REMI02</u>
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q3495</u>
Instrument ID:	<u>ECD_D</u>	Calibration Date(s):	<u>10/17/2025</u> <u>10/17/2025</u>
		Calibration Times:	<u>11:34</u> <u>12:28</u>

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

LAB FILE ID:	RT 100 = <u>PD090650.D</u>	RT 075 = <u>PD090651.D</u>
RT 050 = <u>PD090652.D</u>	RT 025 = <u>PD090653.D</u>	RT 005 = <u>PD090654.D</u>

COMPOUND	RT 100	RT 075	RT 050	RT 025	RT 005	MEAN RT	RT WINDOW	
							FROM	TO
4,4'-DDD	5.92	5.92	5.92	5.92	5.92	5.92	5.82	6.02
4,4'-DDE	5.37	5.37	5.37	5.37	5.37	5.37	5.27	5.47
4,4'-DDT	6.17	6.18	6.17	6.17	6.18	6.17	6.07	6.27
Aldrin	4.36	4.36	4.36	4.36	4.36	4.36	4.26	4.46
alpha-BHC	3.39	3.39	3.39	3.39	3.39	3.39	3.29	3.49
alpha-Chlordane	5.18	5.18	5.18	5.18	5.18	5.18	5.08	5.28
beta-BHC	4.02	4.02	4.02	4.02	4.02	4.02	3.92	4.12
Decachlorobiphenyl	8.06	8.06	8.06	8.06	8.06	8.06	7.96	8.16
delta-BHC	4.26	4.26	4.26	4.26	4.26	4.26	4.16	4.36
Dieldrin	5.50	5.50	5.50	5.50	5.50	5.50	5.40	5.60
Endosulfan I	5.24	5.24	5.24	5.24	5.24	5.24	5.14	5.34
Endosulfan II	6.07	6.07	6.07	6.07	6.07	6.07	5.97	6.17
Endosulfan sulfate	6.47	6.47	6.47	6.47	6.47	6.47	6.37	6.57
Endrin	5.78	5.78	5.78	5.78	5.78	5.78	5.68	5.88
Endrin aldehyde	6.25	6.25	6.25	6.25	6.25	6.25	6.15	6.35
Endrin ketone	6.98	6.98	6.98	6.98	6.98	6.98	6.88	7.08
gamma-BHC (Lindane)	3.72	3.72	3.72	3.72	3.72	3.72	3.62	3.82
gamma-Chlordane	5.12	5.12	5.12	5.12	5.12	5.12	5.02	5.22
Heptachlor	4.08	4.08	4.08	4.07	4.08	4.07	3.97	4.17
Heptachlor epoxide	4.86	4.87	4.86	4.86	4.86	4.86	4.76	4.96
Methoxychlor	6.75	6.75	6.75	6.74	6.75	6.74	6.64	6.84
Tetrachloro-m-xylene	2.87	2.88	2.88	2.87	2.87	2.87	2.77	2.97

CALIBRATION FACTOR OF INITIAL CALIBRATION

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

Contract: REMI02
SDG NO.: Q3495

Calibration Date(s): 10/17/2025 10/17/2025
Calibration Times: 11:34 12:28

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

LAB FILE ID:		CF 100 =	<u>PD090650.D</u>	CF 075 =	<u>PD090651.D</u>		
CF 050 =		<u>PD090652.D</u>	CF 025 =	<u>PD090653.D</u>	CF 005 =	<u>PD090654.D</u>	
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	4150980000	4075270000	4055470000	3840760000	4148570000	4054210000	3
4,4'-DDE	5435610000	5370750000	5317610000	5053060000	5456160000	5326640000	3
4,4'-DDT	4256810000	4152600000	4117930000	3927880000	4097130000	4110470000	3
Aldrin	6694230000	6562230000	6516940000	6238140000	6724820000	6547270000	3
alpha-BHC	7752530000	7402380000	7252210000	6743070000	6735520000	7177140000	6
alpha-Chlordane	5966070000	5905470000	5938650000	6170460000	6920630000	6180250000	7
beta-BHC	2515210000	2489310000	2540370000	2560620000	2992650000	2619630000	8
Decachlorobiphenyl	4361180000	4315270000	4506230000	4707060000	5500070000	4677960000	10
delta-BHC	7169260000	6905720000	6771510000	6342450000	6619680000	6761720000	5
Dieldrin	5929890000	5841940000	5818180000	5550330000	6016400000	5831350000	3
Endosulfan I	5452270000	5404490000	5449050000	5348660000	6028770000	5536650000	5
Endosulfan II	4873230000	4835950000	4885080000	4789750000	5615800000	4999960000	7
Endosulfan sulfate	4557060000	4504850000	4556170000	4479960000	5092920000	4638190000	6
Endrin	5035940000	4962910000	4938250000	4711180000	5120260000	4953710000	3
Endrin aldehyde	3505880000	3507960000	3571680000	3579710000	4148280000	3662700000	7
Endrin ketone	4819770000	4736510000	4782320000	4695250000	5260400000	4858850000	5
gamma-BHC (Lindane)	7185880000	6905860000	6799650000	6417980000	6684750000	6798820000	4
gamma-Chlordane	5972450000	5894040000	5855640000	5592670000	6341090000	5931180000	5
Heptachlor	5757910000	5597510000	5545540000	5288860000	5757970000	5589560000	3
Heptachlor epoxide	5795570000	5708170000	5718680000	5584980000	6137920000	5789060000	4
Methoxychlor	2050060000	2015340000	2078640000	2094570000	2389310000	2125590000	7
Tetrachloro-m-xylene	3010860000	2951180000	3030100000	3124300000	3454390000	3114160000	6

CALIBRATION FACTOR OF INITIAL CALIBRATION

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

Contract: REMI02
SDG NO.: Q3495

Calibration Date(s): 10/17/2025 10/17/2025
Calibration Times: 11:34 12:28

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

LAB FILE ID:		CF 100 =	<u>PD090650.D</u>	CF 075 =	<u>PD090651.D</u>		
CF 050 =		<u>PD090652.D</u>	CF 025 =	<u>PD090653.D</u>	CF 005 =	<u>PD090654.D</u>	
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	33708900000	34309400000	35560900000	36237500000	42373300000	36438000000	10
4,4'-DDE	40500800000	41191000000	42548600000	43577200000	51373700000	43838300000	10
4,4'-DDT	33888000000	33982600000	34868200000	34959100000	37667600000	35073100000	4
Aldrin	43083800000	43574800000	45033800000	46172400000	53729700000	46318900000	9
alpha-BHC	49118000000	48392000000	49683700000	50500300000	57784900000	51095800000	7
alpha-Chlordane	40004000000	40550200000	41893000000	42708900000	50532500000	43137700000	10
beta-BHC	18296700000	18365500000	19080800000	19804100000	23600800000	19829500000	11
Decachlorobiphenyl	28856800000	28247200000	29308000000	30057200000	34987600000	30291400000	9
delta-BHC	45324900000	45220200000	46466400000	47155800000	54416000000	47716700000	8
Dieldrin	40601100000	41425500000	43096200000	44203200000	51593900000	44184000000	10
Endosulfan I	35728600000	36586400000	38185700000	39478600000	46972400000	39390300000	11
Endosulfan II	34269700000	34968100000	36432200000	37520200000	44587200000	37555500000	11
Endosulfan sulfate	32963400000	33538900000	34960800000	35861000000	41900900000	35845000000	10
Endrin	36902000000	37827400000	39400900000	40693000000	47646500000	40494000000	11
Endrin aldehyde	25110900000	25739400000	26893100000	27767400000	32953200000	27692800000	11
Endrin ketone	34460000000	35020500000	36550700000	37816700000	44740900000	37717800000	11
gamma-BHC (Lindane)	44964800000	44604900000	45972800000	46808300000	53893500000	47248900000	8
gamma-Chlordane	41920800000	42422300000	43748100000	44346600000	52306700000	44948900000	9
Heptachlor	41272300000	41810600000	43239300000	44502900000	52189000000	44602800000	10
Heptachlor epoxide	38203400000	38891500000	40672900000	42374400000	53052500000	42639000000	14
Methoxychlor	16116900000	16316400000	17198900000	17735100000	19771700000	17427800000	8
Tetrachloro-m-xylene	27903400000	27689100000	28617400000	29726500000	34949400000	29777100000	10

INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Lab Name: Alliance
Contract: REMI02

Lab Code: ACE
SDG NO.: Q3495

Instrument ID: ECD_D
Date(s) Analyzed: 10/17/2025 10/17/2025

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	32564400
		2	6.44	6.34	6.54	48621800
		3	7.14	7.04	7.24	85710600
		4	7.56	7.46	7.66	95737400
		5	7.92	7.82	8.02	60480500

INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Lab Name: Alliance
Contract: REMI02

Lab Code: ACE
SDG NO.: Q3495

Instrument ID: ECD_D
Date(s) Analyzed: 10/17/2025 10/17/2025

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	263702000
		2	5.64	5.54	5.74	176560000
		3	6.75	6.65	6.85	658883000
		4	7.19	7.09	7.29	495104000
		5	7.32	7.22	7.42	246033000



Lab Name:	<u>Alliance</u>	Contract:	<u>REMI02</u>	
Lab Code:	<u>ACE</u>	SDG NO.:	<u>Q3495</u>	
Instrument ID:	<u>ECD_L</u>	Calibration Date(s):	<u>10/28/2025</u>	<u>10/28/2025</u>
		Calibration Times:	<u>14:57</u>	<u>15:52</u>

LAB FILE ID:	RT 100 = <u>PL097639.D</u>	RT 075 = <u>PL097640.D</u>
RT 050 = <u>PL097641.D</u>	RT 025 = <u>PL097642.D</u>	RT 005 = <u>PL097643.D</u>

[illegible]

A
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L

Contract: REMI02

SDG NO.: Q3495

Calibration Date(s): 10/28/2025 10/28/2025

Calibration Times: 14:57 15:52

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

LAB FILE ID:		RT 100 =	<u>PL097639.D</u>	RT 075 =	<u>PL097640.D</u>
RT 050 =	PL097641.D	RT 025 =	PL097642.D	RT 005 =	PL097643.D

[illegible]

CALIBRATION FACTOR OF INITIAL CALIBRATION

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

Contract: REMI02
SDG NO.: Q3495

Calibration Date(s): 10/28/2025 10/28/2025
Calibration Times: 14:57 15:52

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

LAB FILE ID:		CF 100 =	<u>PL097639.D</u>	CF 075 =	<u>PL097640.D</u>		
CF 050 =		<u>PL097641.D</u>	CF 025 =	<u>PL097642.D</u>	CF 005 =	<u>PL097643.D</u>	
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	3567700000	3596420000	3519300000	3523300000	3666310000	3574610000	2
4,4'-DDE	4662610000	4685370000	4441220000	4360150000	4797990000	4589470000	4
4,4'-DDT	4001630000	4034700000	3952490000	3902070000	4268650000	4031910000	4
Aldrin	6058100000	6091030000	5945310000	5940210000	6490500000	6105030000	4
alpha-BHC	6928520000	6914720000	6685810000	6608510000	6833540000	6794220000	2
alpha-Chlordane	5318710000	5392920000	5297330000	5359280000	6080300000	5489710000	6
beta-BHC	2483420000	2529180000	2519150000	2622160000	2877130000	2606210000	6
Decachlorobiphenyl	2752980000	2829430000	2855460000	2954080000	3483880000	2975170000	10
delta-BHC	5971560000	5952370000	5775720000	5721570000	6304030000	5945050000	4
Dieldrin	5283750000	5330610000	5235600000	5266320000	5691530000	5361560000	4
Endosulfan I	4923010000	4986140000	4874360000	4922540000	5508940000	5043000000	5
Endosulfan II	4327360000	4378110000	4392730000	4521570000	5144310000	4552820000	7
Endosulfan sulfate	3758840000	3821150000	3776740000	3833480000	4357140000	3909470000	6
Endrin	4454260000	4378530000	4274020000	4464660000	4827220000	4479740000	5
Endrin aldehyde	3101220000	3164730000	3160050000	3271010000	3752980000	3290000000	8
Endrin ketone	4059380000	4127980000	4091630000	4143360000	4391020000	4162670000	3
gamma-BHC (Lindane)	6439530000	6439050000	6256040000	6236810000	6572420000	6388770000	2
gamma-Chlordane	5419060000	5469220000	5358210000	5341330000	5823080000	5482180000	4
Heptachlor	6313730000	6354930000	6248720000	6339570000	6718330000	6395060000	3
Heptachlor epoxide	5349750000	5417960000	5398670000	5652150000	7338600000	5831430000	15
Methoxychlor	1916410000	1969950000	1972250000	2019390000	2326100000	2040820000	8
Tetrachloro-m-xylene	3969760000	4019030000	3978350000	4085520000	4406630000	4091860000	4

CALIBRATION FACTOR OF INITIAL CALIBRATION

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

Contract: REMI02
SDG NO.: Q3495

Calibration Date(s): 10/28/2025 10/28/2025
Calibration Times: 14:57 15:52

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

LAB FILE ID:		CF 100 =	<u>PL097639.D</u>	CF 075 =	<u>PL097640.D</u>		
CF 050 =		<u>PL097641.D</u>	CF 025 =	<u>PL097642.D</u>	CF 005 =	<u>PL097643.D</u>	
COMPOUND	CF 100	CF 075	CF 050	CF 025	CF 005	CF	% RSD
4,4'-DDD	6275870000	6447250000	6304230000	6280450000	6746180000	6410790000	3
4,4'-DDE	7606260000	7779780000	7589550000	7495840000	7855120000	7665310000	2
4,4'-DDT	6819710000	6969130000	6786060000	6778330000	7038460000	6878340000	2
Aldrin	8434770000	8618140000	8473930000	8459580000	8998980000	8597080000	3
alpha-BHC	9819830000	9929670000	9677710000	9562720000	9654590000	9728900000	1
alpha-Chlordane	7955180000	8182630000	8114990000	8327570000	8936440000	8303360000	5
beta-BHC	3699680000	3793400000	3790740000	3929010000	4543610000	3951290000	9
Decachlorobiphenyl	4847190000	5011950000	5033010000	5379370000	6452460000	5344800000	12
delta-BHC	9007400000	9145200000	8903490000	8797930000	9020540000	8974910000	1
Dieldrin	7848960000	8071420000	7933490000	7923410000	8420040000	8039460000	3
Endosulfan I	7063640000	7307870000	7289790000	7533680000	8447240000	7528440000	7
Endosulfan II	6655110000	6864970000	6820360000	6950360000	7820390000	7022240000	7
Endosulfan sulfate	6329550000	6567420000	6538450000	6724860000	7571530000	6746360000	7
Endrin	6696330000	6920810000	6835150000	6953450000	7577070000	6996560000	5
Endrin aldehyde	4975120000	5183660000	5236320000	5611370000	7771040000	5755500000	20
Endrin ketone	6941110000	7218950000	7210830000	7499260000	8349000000	7443830000	7
gamma-BHC (Lindane)	9025580000	9156440000	8952920000	8943570000	9216610000	9059020000	1
gamma-Chlordane	8266210000	8501510000	8457090000	8767920000	9279040000	8654350000	5
Heptachlor	8675510000	8892850000	8832330000	8964890000	9961270000	9065370000	6
Heptachlor epoxide	7443930000	7693490000	7641160000	7770100000	8710360000	7851810000	6
Methoxychlor	3414390000	3555870000	3587550000	3765390000	4253070000	3715250000	9
Tetrachloro-m-xylene	5762550000	5872210000	5827140000	5996140000	6637090000	6019020000	6

INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Instrument ID: ECD_L

Date(s) Analyzed: 10/28/2025 10/28/2025

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

COMPOUND	AMOUNT (ng)	PEAK	RT	RT WINDOW		CALIBRATION FACTOR
				FROM	TO	
Toxaphene	500	1	6.20	6.10	6.30	40325100
		2	6.60	6.50	6.70	29864600
		3	7.01	6.91	7.11	141530000
		4	7.10	7.00	7.20	102524000
		5	7.88	7.78	7.98	73398900

INITIAL CALIBRATION OF MULTICOMPONENT ANALYTES

Lab Name: Alliance Contract: REMI02

Lab Code: ACE SDG NO.: Q3495

Instrument ID: ECD_L Date(s) Analyzed: 10/28/2025 10/28/2025

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

COMPOUND	AMOUNT (ng)	PEAK	RT	RT WINDOW		CALIBRATION FACTOR
				FROM	TO	
Toxaphene	500	1	5.06	4.96	5.16	42932600
		2	5.75	5.65	5.85	48959900
		3	6.03	5.93	6.13	56151700
		4	6.66	6.56	6.76	172906000
		5	7.10	7.00	7.20	107207000

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/17/2025 10/17/2025

Continuing Calib Time: 11:09

Initial Calibration Time(s): 11:34 12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.08	9.06	8.96	9.16	-0.02
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	3.99	3.89	4.09	-0.01
beta-BHC	4.52	4.51	4.41	4.61	-0.01
delta-BHC	4.77	4.76	4.66	4.86	-0.01
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.92	4.82	5.02	-0.01
Aldrin	5.27	5.26	5.16	5.36	-0.01
Heptachlor epoxide	5.70	5.68	5.58	5.78	-0.02
Endosulfan I	6.08	6.07	5.97	6.17	-0.01
Dieldrin	6.35	6.34	6.24	6.44	-0.01
4,4'-DDE	6.20	6.19	6.09	6.29	-0.01
Endrin	6.58	6.57	6.47	6.67	-0.01
Endosulfan II	6.79	6.78	6.68	6.88	-0.01
4,4'-DDD	6.71	6.70	6.60	6.80	-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	6.91	6.81	7.01	-0.01
alpha-Chlordane	6.03	6.02	5.92	6.12	-0.01
gamma-Chlordane	5.95	5.94	5.84	6.04	-0.01

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/17/2025 10/17/2025

Continuing Calib Time: 11:09

Initial Calibration Time(s): 11:34 12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.07	8.06	7.96	8.16	-0.01
Tetrachloro-m-xylene	2.87	2.88	2.78	2.98	0.01
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.02	4.02	3.92	4.12	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.72	3.72	3.62	3.82	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.01
Aldrin	4.36	4.36	4.26	4.46	0.00
Heptachlor epoxide	4.87	4.86	4.76	4.96	0.00
Endosulfan I	5.24	5.24	5.14	5.34	0.00
Dieldrin	5.51	5.50	5.40	5.60	-0.01
4,4'-DDE	5.37	5.37	5.27	5.47	0.00
Endrin	5.78	5.78	5.68	5.88	0.00
Endosulfan II	6.07	6.07	5.97	6.17	0.00
4,4'-DDD	5.92	5.92	5.82	6.02	0.00
Endosulfan sulfate	6.48	6.47	6.37	6.57	-0.01
4,4'-DDT	6.18	6.17	6.07	6.27	-0.01
Methoxychlor	6.75	6.75	6.65	6.85	0.00
Endrin ketone	6.99	6.98	6.88	7.08	-0.01
Endrin aldehyde	6.25	6.25	6.15	6.35	0.00
alpha-Chlordane	5.18	5.18	5.08	5.28	0.00
gamma-Chlordane	5.12	5.12	5.02	5.22	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL01 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090926.D **Time Analyzed:** 11:09

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.712	6.600	6.800	48.540	50.000	-2.9
4,4'-DDE	6.202	6.090	6.290	45.040	50.000	-9.9
4,4'-DDT	7.028	6.915	7.115	51.870	50.000	3.7
Aldrin	5.274	5.164	5.364	45.260	50.000	-9.5
alpha-BHC	4.003	3.894	4.094	45.180	50.000	-9.6
alpha-Chlordane	6.032	5.921	6.121	43.400	50.000	-13.2
beta-BHC	4.522	4.412	4.612	43.330	50.000	-13.3
Decachlorobiphenyl	9.081	8.964	9.164	43.690	50.000	-12.6
delta-BHC	4.770	4.660	4.860	45.000	50.000	-10.0
Dieldrin	6.352	6.240	6.440	46.130	50.000	-7.7
Endosulfan I	6.079	5.968	6.168	45.300	50.000	-9.4
Endosulfan II	6.792	6.681	6.881	46.350	50.000	-7.3
Endosulfan sulfate	7.156	7.045	7.245	45.990	50.000	-8.0
Endrin	6.580	6.468	6.668	46.810	50.000	-6.4
Endrin aldehyde	6.922	6.810	7.010	47.350	50.000	-5.3
Endrin ketone	7.637	7.525	7.725	47.160	50.000	-5.7
gamma-BHC (Lindane)	4.334	4.225	4.425	44.400	50.000	-11.2
gamma-Chlordane	5.951	5.840	6.040	45.120	50.000	-9.8
Heptachlor	4.932	4.823	5.023	56.000	50.000	12.0
Heptachlor epoxide	5.695	5.584	5.784	44.950	50.000	-10.1
Methoxychlor	7.501	7.388	7.588	52.280	50.000	4.6
Tetrachloro-m-xylene	3.552	3.445	3.645	42.890	50.000	-14.2

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL01 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090926.D **Time Analyzed:** 11:09

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.923	5.821	6.021	50.860	50.000	1.7
4,4'-DDE	5.367	5.266	5.466	50.270	50.000	0.5
4,4'-DDT	6.177	6.074	6.274	57.220	50.000	14.4
Aldrin	4.361	4.261	4.461	49.550	50.000	-0.9
alpha-BHC	3.386	3.287	3.487	49.450	50.000	-1.1
alpha-Chlordane	5.183	5.081	5.281	50.040	50.000	0.1
beta-BHC	4.019	3.919	4.119	48.710	50.000	-2.6
Decachlorobiphenyl	8.066	7.962	8.162	56.820	50.000	13.6
delta-BHC	4.255	4.155	4.355	49.160	50.000	-1.7
Dieldrin	5.506	5.404	5.604	50.670	50.000	1.3
Endosulfan I	5.240	5.138	5.338	48.720	50.000	-2.6
Endosulfan II	6.074	5.972	6.172	52.200	50.000	4.4
Endosulfan sulfate	6.477	6.374	6.574	52.480	50.000	5.0
Endrin	5.783	5.681	5.881	51.560	50.000	3.1
Endrin aldehyde	6.253	6.150	6.350	52.780	50.000	5.6
Endrin ketone	6.985	6.883	7.083	58.590	50.000	17.2
gamma-BHC (Lindane)	3.723	3.623	3.823	49.340	50.000	-1.3
gamma-Chlordane	5.118	5.017	5.217	50.120	50.000	0.2
Heptachlor	4.075	3.975	4.175	51.770	50.000	3.5
Heptachlor epoxide	4.865	4.764	4.964	48.690	50.000	-2.6
Methoxychlor	6.748	6.645	6.845	56.010	50.000	12.0
Tetrachloro-m-xylene	2.873	2.775	2.975	48.290	50.000	-3.4

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/17/2025

10/17/2025

Continuing Calib Time: 16:01

Initial Calibration Time(s): 11:34

12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.06	8.96	9.16	-0.01
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	3.99	3.89	4.09	0.00
beta-BHC	4.51	4.51	4.41	4.61	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	4.92	4.92	4.82	5.02	0.00
Aldrin	5.27	5.26	5.16	5.36	0.00
Heptachlor epoxide	5.69	5.68	5.58	5.78	-0.01
Endosulfan I	6.07	6.07	5.97	6.17	0.00
Dieldrin	6.34	6.34	6.24	6.44	0.00
4,4'-DDE	6.19	6.19	6.09	6.29	0.00
Endrin	6.57	6.57	6.47	6.67	0.00
Endosulfan II	6.78	6.78	6.68	6.88	0.00
4,4'-DDD	6.70	6.70	6.60	6.80	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	6.91	6.91	6.81	7.01	0.00
alpha-Chlordane	6.02	6.02	5.92	6.12	0.00
gamma-Chlordane	5.94	5.94	5.84	6.04	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/17/2025

10/17/2025

Continuing Calib Time: 16:01

Initial Calibration Time(s): 11:34

12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.06	8.06	7.96	8.16	0.00
Tetrachloro-m-xylene	2.87	2.88	2.78	2.98	0.01
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.02	4.02	3.92	4.12	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.72	3.72	3.62	3.82	0.00
Heptachlor	4.08	4.08	3.98	4.18	0.01
Aldrin	4.36	4.36	4.26	4.46	0.00
Heptachlor epoxide	4.86	4.86	4.76	4.96	0.00
Endosulfan I	5.24	5.24	5.14	5.34	0.00
Dieldrin	5.51	5.50	5.40	5.60	-0.01
4,4'-DDE	5.37	5.37	5.27	5.47	0.00
Endrin	5.78	5.78	5.68	5.88	0.00
Endosulfan II	6.07	6.07	5.97	6.17	0.00
4,4'-DDD	5.92	5.92	5.82	6.02	0.00
Endosulfan sulfate	6.47	6.47	6.37	6.57	0.00
4,4'-DDT	6.17	6.17	6.07	6.27	0.00
Methoxychlor	6.75	6.75	6.65	6.85	0.01
Endrin ketone	6.98	6.98	6.88	7.08	0.00
Endrin aldehyde	6.25	6.25	6.15	6.35	0.00
alpha-Chlordane	5.18	5.18	5.08	5.28	0.00
gamma-Chlordane	5.12	5.12	5.02	5.22	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL02 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090933.D **Time Analyzed:** 16:01

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.702	6.600	6.800	46.530	50.000	-6.9
4,4'-DDE	6.192	6.090	6.290	43.680	50.000	-12.6
4,4'-DDT	7.018	6.915	7.115	50.220	50.000	0.4
Aldrin	5.265	5.164	5.364	44.350	50.000	-11.3
alpha-BHC	3.995	3.894	4.094	44.810	50.000	-10.4
alpha-Chlordane	6.022	5.921	6.121	42.650	50.000	-14.7
beta-BHC	4.513	4.412	4.612	43.330	50.000	-13.3
Decachlorobiphenyl	9.070	8.964	9.164	42.870	50.000	-14.3
delta-BHC	4.761	4.660	4.860	43.970	50.000	-12.1
Dieldrin	6.343	6.240	6.440	45.000	50.000	-10.0
Endosulfan I	6.070	5.968	6.168	44.180	50.000	-11.6
Endosulfan II	6.783	6.681	6.881	44.920	50.000	-10.2
Endosulfan sulfate	7.147	7.045	7.245	44.700	50.000	-10.6
Endrin	6.570	6.468	6.668	45.550	50.000	-8.9
Endrin aldehyde	6.912	6.810	7.010	46.010	50.000	-8.0
Endrin ketone	7.628	7.525	7.725	45.830	50.000	-8.3
gamma-BHC (Lindane)	4.326	4.225	4.425	44.280	50.000	-11.4
gamma-Chlordane	5.942	5.840	6.040	43.940	50.000	-12.1
Heptachlor	4.924	4.823	5.023	54.600	50.000	9.2
Heptachlor epoxide	5.686	5.584	5.784	44.780	50.000	-10.4
Methoxychlor	7.491	7.388	7.588	50.720	50.000	1.4
Tetrachloro-m-xylene	3.545	3.445	3.645	43.080	50.000	-13.8

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL02 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090933.D **Time Analyzed:** 16:01

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.921	5.821	6.021	50.070	50.000	0.1
4,4'-DDE	5.366	5.266	5.466	49.650	50.000	-0.7
4,4'-DDT	6.174	6.074	6.274	56.020	50.000	12.0
Aldrin	4.360	4.261	4.461	49.190	50.000	-1.6
alpha-BHC	3.386	3.287	3.487	48.830	50.000	-2.3
alpha-Chlordane	5.181	5.081	5.281	49.170	50.000	-1.7
beta-BHC	4.019	3.919	4.119	48.210	50.000	-3.6
Decachlorobiphenyl	8.062	7.962	8.162	56.070	50.000	12.1
delta-BHC	4.255	4.155	4.355	48.750	50.000	-2.5
Dieldrin	5.505	5.404	5.604	50.000	50.000	0.0
Endosulfan I	5.238	5.138	5.338	47.980	50.000	-4.0
Endosulfan II	6.072	5.972	6.172	50.960	50.000	1.9
Endosulfan sulfate	6.474	6.374	6.574	51.370	50.000	2.7
Endrin	5.781	5.681	5.881	51.480	50.000	3.0
Endrin aldehyde	6.251	6.150	6.350	51.790	50.000	3.6
Endrin ketone	6.983	6.883	7.083	56.530	50.000	13.1
gamma-BHC (Lindane)	3.723	3.623	3.823	48.820	50.000	-2.4
gamma-Chlordane	5.117	5.017	5.217	49.240	50.000	-1.5
Heptachlor	4.075	3.975	4.175	51.650	50.000	3.3
Heptachlor epoxide	4.864	4.764	4.964	48.020	50.000	-4.0
Methoxychlor	6.745	6.645	6.845	56.430	50.000	12.9
Tetrachloro-m-xylene	2.874	2.775	2.975	47.690	50.000	-4.6

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 11/04/2025

Initial Calibration Date(s): 10/17/2025 10/17/2025

Continuing Calib Time: 13:29

Initial Calibration Time(s): 11:34 12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.07	9.06	8.96	9.16	-0.01
Tetrachloro-m-xylene	3.55	3.55	3.45	3.65	0.00
alpha-BHC	4.00	3.99	3.89	4.09	-0.01
beta-BHC	4.52	4.51	4.41	4.61	-0.01
delta-BHC	4.77	4.76	4.66	4.86	-0.01
gamma-BHC (Lindane)	4.33	4.33	4.23	4.43	0.00
Heptachlor	4.93	4.92	4.82	5.02	-0.01
Aldrin	5.27	5.26	5.16	5.36	-0.01
Heptachlor epoxide	5.69	5.68	5.58	5.78	-0.01
Endosulfan I	6.07	6.07	5.97	6.17	0.00
Dieldrin	6.35	6.34	6.24	6.44	-0.01
4,4'-DDE	6.20	6.19	6.09	6.29	-0.01
Endrin	6.58	6.57	6.47	6.67	0.00
Endosulfan II	6.79	6.78	6.68	6.88	-0.01
4,4'-DDD	6.71	6.70	6.60	6.80	-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	7.50	7.49	7.39	7.59	0.00
Endrin ketone	7.63	7.63	7.53	7.73	0.00
Endrin aldehyde	6.92	6.91	6.81	7.01	-0.01
alpha-Chlordane	6.03	6.02	5.92	6.12	-0.01
gamma-Chlordane	5.95	5.94	5.84	6.04	-0.01

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 11/04/2025

Initial Calibration Date(s): 10/17/2025

10/17/2025

Continuing Calib Time: 13:29

Initial Calibration Time(s): 11:34

12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.06	8.06	7.96	8.16	0.00
Tetrachloro-m-xylene	2.87	2.88	2.78	2.98	0.01
alpha-BHC	3.39	3.39	3.29	3.49	0.00
beta-BHC	4.02	4.02	3.92	4.12	0.00
delta-BHC	4.26	4.26	4.16	4.36	0.00
gamma-BHC (Lindane)	3.72	3.72	3.62	3.82	0.00
Heptachlor	4.07	4.08	3.98	4.18	0.01
Aldrin	4.36	4.36	4.26	4.46	0.00
Heptachlor epoxide	4.86	4.86	4.76	4.96	0.00
Endosulfan I	5.24	5.24	5.14	5.34	0.00
Dieldrin	5.50	5.50	5.40	5.60	0.00
4,4'-DDE	5.37	5.37	5.27	5.47	0.01
Endrin	5.78	5.78	5.68	5.88	0.00
Endosulfan II	6.07	6.07	5.97	6.17	0.00
4,4'-DDD	5.92	5.92	5.82	6.02	0.00
Endosulfan sulfate	6.47	6.47	6.37	6.57	0.00
4,4'-DDT	6.17	6.17	6.07	6.27	0.00
Methoxychlor	6.75	6.75	6.65	6.85	0.01
Endrin ketone	6.98	6.98	6.88	7.08	0.00
Endrin aldehyde	6.25	6.25	6.15	6.35	0.00
alpha-Chlordane	5.18	5.18	5.08	5.28	0.00
gamma-Chlordane	5.12	5.12	5.02	5.22	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL03 **Date Analyzed:** 11/04/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090949.D **Time Analyzed:** 13:29

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.706	6.600	6.800	51.660	50.000	3.3
4,4'-DDE	6.197	6.090	6.290	48.680	50.000	-2.6
4,4'-DDT	7.022	6.915	7.115	48.520	50.000	-3.0
Aldrin	5.270	5.164	5.364	52.900	50.000	5.8
alpha-BHC	4.000	3.894	4.094	55.710	50.000	11.4
alpha-Chlordane	6.027	5.921	6.121	47.950	50.000	-4.1
beta-BHC	4.519	4.412	4.612	51.550	50.000	3.1
Decachlorobiphenyl	9.074	8.964	9.164	43.990	50.000	-12.0
delta-BHC	4.767	4.660	4.860	54.270	50.000	8.5
Dieldrin	6.347	6.240	6.440	50.090	50.000	0.2
Endosulfan I	6.074	5.968	6.168	49.820	50.000	-0.4
Endosulfan II	6.787	6.681	6.881	48.630	50.000	-2.7
Endosulfan sulfate	7.151	7.045	7.245	45.780	50.000	-8.4
Endrin	6.575	6.468	6.668	49.110	50.000	-1.8
Endrin aldehyde	6.917	6.810	7.010	45.580	50.000	-8.8
Endrin ketone	7.631	7.525	7.725	47.380	50.000	-5.2
gamma-BHC (Lindane)	4.332	4.225	4.425	54.710	50.000	9.4
gamma-Chlordane	5.947	5.840	6.040	50.180	50.000	0.4
Heptachlor	4.929	4.823	5.023	66.730	50.000	33.5
Heptachlor epoxide	5.690	5.584	5.784	49.950	50.000	-0.1
Methoxychlor	7.495	7.388	7.588	50.020	50.000	0.0
Tetrachloro-m-xylene	3.551	3.445	3.645	51.620	50.000	3.2

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL03 **Date Analyzed:** 11/04/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090949.D **Time Analyzed:** 13:29

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.920	5.821	6.021	47.850	50.000	-4.3
4,4'-DDE	5.365	5.266	5.466	49.430	50.000	-1.1
4,4'-DDT	6.174	6.074	6.274	48.230	50.000	-3.5
Aldrin	4.359	4.261	4.461	52.620	50.000	5.2
alpha-BHC	3.386	3.287	3.487	53.750	50.000	7.5
alpha-Chlordane	5.181	5.081	5.281	50.060	50.000	0.1
beta-BHC	4.018	3.919	4.119	52.130	50.000	4.3
Decachlorobiphenyl	8.062	7.962	8.162	37.880	50.000	-24.2
delta-BHC	4.255	4.155	4.355	52.950	50.000	5.9
Dieldrin	5.503	5.404	5.604	49.200	50.000	-1.6
Endosulfan I	5.238	5.138	5.338	48.710	50.000	-2.6
Endosulfan II	6.072	5.972	6.172	47.370	50.000	-5.3
Endosulfan sulfate	6.474	6.374	6.574	43.940	50.000	-12.1
Endrin	5.780	5.681	5.881	47.270	50.000	-5.5
Endrin aldehyde	6.250	6.150	6.350	44.960	50.000	-10.1
Endrin ketone	6.983	6.883	7.083	40.460	50.000	-19.1
gamma-BHC (Lindane)	3.722	3.623	3.823	53.140	50.000	6.3
gamma-Chlordane	5.116	5.017	5.217	50.480	50.000	1.0
Heptachlor	4.074	3.975	4.175	55.600	50.000	11.2
Heptachlor epoxide	4.864	4.764	4.964	51.470	50.000	2.9
Methoxychlor	6.745	6.645	6.845	43.330	50.000	-13.3
Tetrachloro-m-xylene	2.874	2.775	2.975	53.000	50.000	6.0

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 11/04/2025

Initial Calibration Date(s): 10/17/2025

10/17/2025

Continuing Calib Time: 16:44

Initial Calibration Time(s): 11:34

12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.06	9.06	8.96	9.16	0.00
Tetrachloro-m-xylene	3.54	3.55	3.45	3.65	0.01
alpha-BHC	3.99	3.99	3.89	4.09	0.00
beta-BHC	4.51	4.51	4.41	4.61	0.00
delta-BHC	4.76	4.76	4.66	4.86	0.00
gamma-BHC (Lindane)	4.32	4.33	4.23	4.43	0.01
Heptachlor	4.92	4.92	4.82	5.02	0.00
Aldrin	5.26	5.26	5.16	5.36	0.00
Heptachlor epoxide	5.68	5.68	5.58	5.78	0.00
Endosulfan I	6.07	6.07	5.97	6.17	0.00
Dieldrin	6.34	6.34	6.24	6.44	0.00
4,4'-DDE	6.19	6.19	6.09	6.29	0.00
Endrin	6.57	6.57	6.47	6.67	0.00
Endosulfan II	6.78	6.78	6.68	6.88	0.00
4,4'-DDD	6.70	6.70	6.60	6.80	0.00
Endosulfan sulfate	7.14	7.15	7.05	7.25	0.01
4,4'-DDT	7.02	7.02	6.92	7.12	0.01
Methoxychlor	7.49	7.49	7.39	7.59	0.00
Endrin ketone	7.62	7.63	7.53	7.73	0.01
Endrin aldehyde	6.91	6.91	6.81	7.01	0.00
alpha-Chlordane	6.02	6.02	5.92	6.12	0.00
gamma-Chlordane	5.94	5.94	5.84	6.04	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 11/04/2025

Initial Calibration Date(s): 10/17/2025

10/17/2025

Continuing Calib Time: 16:44

Initial Calibration Time(s): 11:34

12:28

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	8.06	8.06	7.96	8.16	0.00
Tetrachloro-m-xylene	2.87	2.88	2.78	2.98	0.01
alpha-BHC	3.39	3.39	3.29	3.49	0.01
beta-BHC	4.02	4.02	3.92	4.12	0.00
delta-BHC	4.25	4.26	4.16	4.36	0.01
gamma-BHC (Lindane)	3.72	3.72	3.62	3.82	0.00
Heptachlor	4.07	4.08	3.98	4.18	0.01
Aldrin	4.36	4.36	4.26	4.46	0.00
Heptachlor epoxide	4.86	4.86	4.76	4.96	0.00
Endosulfan I	5.24	5.24	5.14	5.34	0.00
Dieldrin	5.50	5.50	5.40	5.60	0.00
4,4'-DDE	5.36	5.37	5.27	5.47	0.01
Endrin	5.78	5.78	5.68	5.88	0.00
Endosulfan II	6.07	6.07	5.97	6.17	0.00
4,4'-DDD	5.92	5.92	5.82	6.02	0.00
Endosulfan sulfate	6.47	6.47	6.37	6.57	0.00
4,4'-DDT	6.17	6.17	6.07	6.27	0.00
Methoxychlor	6.74	6.75	6.65	6.85	0.01
Endrin ketone	6.98	6.98	6.88	7.08	0.00
Endrin aldehyde	6.25	6.25	6.15	6.35	0.00
alpha-Chlordane	5.18	5.18	5.08	5.28	0.00
gamma-Chlordane	5.12	5.12	5.02	5.22	0.01

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL04 **Date Analyzed:** 11/04/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090957.D **Time Analyzed:** 16:44

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.699	6.600	6.800	56.440	50.000	12.9
4,4'-DDE	6.189	6.090	6.290	53.890	50.000	7.8
4,4'-DDT	7.015	6.915	7.115	53.170	50.000	6.3
Aldrin	5.263	5.164	5.364	54.610	50.000	9.2
alpha-BHC	3.993	3.894	4.094	55.760	50.000	11.5
alpha-Chlordane	6.020	5.921	6.121	52.830	50.000	5.7
beta-BHC	4.511	4.412	4.612	53.250	50.000	6.5
Decachlorobiphenyl	9.064	8.964	9.164	48.210	50.000	-3.6
delta-BHC	4.760	4.660	4.860	55.270	50.000	10.5
Dieldrin	6.340	6.240	6.440	55.040	50.000	10.1
Endosulfan I	6.067	5.968	6.168	54.280	50.000	8.6
Endosulfan II	6.780	6.681	6.881	52.890	50.000	5.8
Endosulfan sulfate	7.143	7.045	7.245	51.440	50.000	2.9
Endrin	6.567	6.468	6.668	54.530	50.000	9.1
Endrin aldehyde	6.909	6.810	7.010	51.240	50.000	2.5
Endrin ketone	7.624	7.525	7.725	51.440	50.000	2.9
gamma-BHC (Lindane)	4.324	4.225	4.425	55.080	50.000	10.2
gamma-Chlordane	5.939	5.840	6.040	53.960	50.000	7.9
Heptachlor	4.922	4.823	5.023	66.530	50.000	33.1
Heptachlor epoxide	5.683	5.584	5.784	54.960	50.000	9.9
Methoxychlor	7.487	7.388	7.588	52.810	50.000	5.6
Tetrachloro-m-xylene	3.544	3.445	3.645	52.920	50.000	5.8

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/17/2025 10/17/2025

Client Sample No.: CCAL04 **Date Analyzed:** 11/04/2025
Lab Sample No.: PSTDCCC050 **Data File :** PD090957.D **Time Analyzed:** 16:44

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.918	5.821	6.021	54.300	50.000	8.6
4,4'-DDE	5.364	5.266	5.466	52.910	50.000	5.8
4,4'-DDT	6.172	6.074	6.274	52.000	50.000	4.0
Aldrin	4.359	4.261	4.461	52.640	50.000	5.3
alpha-BHC	3.385	3.287	3.487	53.410	50.000	6.8
alpha-Chlordane	5.180	5.081	5.281	52.490	50.000	5.0
beta-BHC	4.018	3.919	4.119	52.430	50.000	4.9
Decachlorobiphenyl	8.060	7.962	8.162	46.000	50.000	-8.0
delta-BHC	4.254	4.155	4.355	53.330	50.000	6.7
Dieldrin	5.502	5.404	5.604	53.290	50.000	6.6
Endosulfan I	5.236	5.138	5.338	51.070	50.000	2.1
Endosulfan II	6.070	5.972	6.172	52.170	50.000	4.3
Endosulfan sulfate	6.472	6.374	6.574	50.680	50.000	1.4
Endrin	5.778	5.681	5.881	53.220	50.000	6.4
Endrin aldehyde	6.248	6.150	6.350	50.540	50.000	1.1
Endrin ketone	6.980	6.883	7.083	49.200	50.000	-1.6
gamma-BHC (Lindane)	3.722	3.623	3.823	53.100	50.000	6.2
gamma-Chlordane	5.115	5.017	5.217	52.640	50.000	5.3
Heptachlor	4.073	3.975	4.175	55.570	50.000	11.1
Heptachlor epoxide	4.862	4.764	4.964	51.570	50.000	3.1
Methoxychlor	6.742	6.645	6.845	51.400	50.000	2.8
Tetrachloro-m-xylene	2.874	2.775	2.975	52.660	50.000	5.3

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/28/2025 10/28/2025

Continuing Calib Time: 11:02

Initial Calibration Time(s): 14:57 15:52

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.01	8.99	8.89	9.09	-0.02
Tetrachloro-m-xylene	3.52	3.52	3.42	3.62	0.00
alpha-BHC	3.97	3.97	3.87	4.07	0.00
beta-BHC	4.49	4.49	4.39	4.59	0.00
delta-BHC	4.73	4.73	4.63	4.83	0.00
gamma-BHC (Lindane)	4.30	4.30	4.20	4.40	0.00
Heptachlor	4.89	4.89	4.79	4.99	0.00
Aldrin	5.23	5.23	5.13	5.33	0.00
Heptachlor epoxide	5.65	5.65	5.55	5.75	0.00
Endosulfan I	6.03	6.03	5.93	6.13	0.00
Dieldrin	6.31	6.30	6.20	6.40	-0.01
4,4'-DDE	6.16	6.15	6.05	6.25	-0.01
Endrin	6.53	6.53	6.43	6.63	0.00
Endosulfan II	6.75	6.74	6.64	6.84	-0.01
4,4'-DDD	6.67	6.66	6.56	6.76	0.00
Endosulfan sulfate	7.11	7.10	7.00	7.20	-0.01
4,4'-DDT	6.98	6.97	6.87	7.07	-0.01
Methoxychlor	7.46	7.45	7.35	7.55	0.00
Endrin ketone	7.59	7.58	7.48	7.68	-0.01
Endrin aldehyde	6.88	6.87	6.77	6.97	-0.01
alpha-Chlordane	5.99	5.98	5.88	6.08	-0.01
gamma-Chlordane	5.90	5.90	5.80	6.00	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/28/2025 10/28/2025

Continuing Calib Time: 11:02

Initial Calibration Time(s): 14:57 15:52

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	7.98	7.98	7.88	8.08	0.00
Tetrachloro-m-xylene	2.82	2.82	2.72	2.92	0.00
alpha-BHC	3.32	3.33	3.23	3.43	0.01
beta-BHC	3.95	3.95	3.85	4.05	0.00
delta-BHC	4.19	4.19	4.09	4.29	0.01
gamma-BHC (Lindane)	3.66	3.66	3.56	3.76	0.01
Heptachlor	4.00	4.01	3.91	4.11	0.01
Aldrin	4.29	4.29	4.19	4.39	0.00
Heptachlor epoxide	4.79	4.79	4.69	4.89	0.00
Endosulfan I	5.16	5.16	5.06	5.26	0.00
Dieldrin	5.42	5.42	5.32	5.52	0.00
4,4'-DDE	5.29	5.29	5.19	5.39	0.00
Endrin	5.70	5.70	5.60	5.80	0.00
Endosulfan II	5.99	5.99	5.89	6.09	0.00
4,4'-DDD	5.85	5.85	5.75	5.95	0.00
Endosulfan sulfate	6.39	6.39	6.29	6.49	0.00
4,4'-DDT	6.10	6.10	6.00	6.20	0.00
Methoxychlor	6.67	6.67	6.57	6.77	0.00
Endrin ketone	6.90	6.90	6.80	7.00	0.00
Endrin aldehyde	6.17	6.17	6.07	6.27	0.00
alpha-Chlordane	5.10	5.10	5.00	5.20	0.00
gamma-Chlordane	5.04	5.04	4.94	5.14	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/28/2025 10/28/2025

Client Sample No.: CCAL05 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PL097701.D **Time Analyzed:** 11:02

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.665	6.560	6.760	56.430	50.000	12.9
4,4'-DDE	6.156	6.051	6.251	56.430	50.000	12.9
4,4'-DDT	6.981	6.874	7.074	51.580	50.000	3.2
Aldrin	5.230	5.127	5.327	54.640	50.000	9.3
alpha-BHC	3.971	3.870	4.070	55.870	50.000	11.7
alpha-Chlordane	5.985	5.881	6.081	54.860	50.000	9.7
beta-BHC	4.488	4.386	4.586	55.240	50.000	10.5
Decachlorobiphenyl	9.005	8.893	9.093	48.940	50.000	-2.1
delta-BHC	4.734	4.631	4.831	58.620	50.000	17.2
Dieldrin	6.306	6.200	6.400	53.420	50.000	6.8
Endosulfan I	6.033	5.928	6.128	55.820	50.000	11.6
Endosulfan II	6.746	6.641	6.841	51.310	50.000	2.6
Endosulfan sulfate	7.109	7.003	7.203	52.430	50.000	4.9
Endrin	6.531	6.427	6.627	53.140	50.000	6.3
Endrin aldehyde	6.875	6.769	6.969	51.160	50.000	2.3
Endrin ketone	7.589	7.481	7.681	54.080	50.000	8.2
gamma-BHC (Lindane)	4.299	4.198	4.398	55.510	50.000	11.0
gamma-Chlordane	5.904	5.801	6.001	53.610	50.000	7.2
Heptachlor	4.889	4.788	4.988	55.480	50.000	11.0
Heptachlor epoxide	5.651	5.547	5.747	53.350	50.000	6.7
Methoxychlor	7.455	7.347	7.547	48.460	50.000	-3.1
Tetrachloro-m-xylene	3.523	3.424	3.624	55.300	50.000	10.6

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/28/2025 10/28/2025

Client Sample No.: CCAL05 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PL097701.D **Time Analyzed:** 11:02

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.847	5.745	5.945	52.580	50.000	5.2
4,4'-DDE	5.293	5.193	5.393	51.700	50.000	3.4
4,4'-DDT	6.099	5.997	6.197	49.420	50.000	-1.2
Aldrin	4.286	4.187	4.387	52.480	50.000	5.0
alpha-BHC	3.323	3.226	3.426	53.470	50.000	6.9
alpha-Chlordane	5.104	5.004	5.204	50.820	50.000	1.6
beta-BHC	3.952	3.854	4.054	52.170	50.000	4.3
Decachlorobiphenyl	7.984	7.879	8.079	43.900	50.000	-12.2
delta-BHC	4.185	4.086	4.286	52.530	50.000	5.1
Dieldrin	5.423	5.324	5.524	51.980	50.000	4.0
Endosulfan I	5.159	5.059	5.259	49.340	50.000	-1.3
Endosulfan II	5.992	5.890	6.090	52.040	50.000	4.1
Endosulfan sulfate	6.394	6.291	6.491	49.350	50.000	-1.3
Endrin	5.699	5.599	5.799	53.140	50.000	6.3
Endrin aldehyde	6.170	6.068	6.268	45.880	50.000	-8.2
Endrin ketone	6.899	6.796	6.996	51.460	50.000	2.9
gamma-BHC (Lindane)	3.655	3.558	3.758	53.110	50.000	6.2
gamma-Chlordane	5.040	4.940	5.140	49.840	50.000	-0.3
Heptachlor	4.003	3.905	4.105	52.530	50.000	5.1
Heptachlor epoxide	4.788	4.689	4.889	51.620	50.000	3.2
Methoxychlor	6.672	6.569	6.769	46.410	50.000	-7.2
Tetrachloro-m-xylene	2.818	2.721	2.921	52.500	50.000	5.0

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/28/2025 10/28/2025

Continuing Calib Time: 16:03

Initial Calibration Time(s): 14:57 15:52

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	9.00	8.99	8.89	9.09	-0.01
Tetrachloro-m-xylene	3.53	3.52	3.42	3.62	-0.01
alpha-BHC	3.97	3.97	3.87	4.07	0.00
beta-BHC	4.49	4.49	4.39	4.59	0.00
delta-BHC	4.73	4.73	4.63	4.83	0.00
gamma-BHC (Lindane)	4.30	4.30	4.20	4.40	0.00
Heptachlor	4.89	4.89	4.79	4.99	0.00
Aldrin	5.23	5.23	5.13	5.33	0.00
Heptachlor epoxide	5.65	5.65	5.55	5.75	0.00
Endosulfan I	6.03	6.03	5.93	6.13	0.00
Dieldrin	6.30	6.30	6.20	6.40	0.00
4,4'-DDE	6.16	6.15	6.05	6.25	-0.01
Endrin	6.53	6.53	6.43	6.63	0.00
Endosulfan II	6.74	6.74	6.64	6.84	0.00
4,4'-DDD	6.66	6.66	6.56	6.76	0.00
Endosulfan sulfate	7.11	7.10	7.00	7.20	-0.01
4,4'-DDT	6.98	6.97	6.87	7.07	-0.01
Methoxychlor	7.45	7.45	7.35	7.55	0.00
Endrin ketone	7.59	7.58	7.48	7.68	-0.01
Endrin aldehyde	6.87	6.87	6.77	6.97	0.00
alpha-Chlordane	5.98	5.98	5.88	6.08	0.00
gamma-Chlordane	5.90	5.90	5.80	6.00	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Continuing Calib Date: 10/31/2025

Initial Calibration Date(s): 10/28/2025 10/28/2025

Continuing Calib Time: 16:03

Initial Calibration Time(s): 14:57 15:52

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

COMPOUND	CCAL RT	AVG RT	RT WINDOW		DIFF RT
			FROM	TO	
Decachlorobiphenyl	7.98	7.98	7.88	8.08	0.00
Tetrachloro-m-xylene	2.82	2.82	2.72	2.92	0.00
alpha-BHC	3.33	3.33	3.23	3.43	0.00
beta-BHC	3.96	3.95	3.85	4.05	0.00
delta-BHC	4.19	4.19	4.09	4.29	0.00
gamma-BHC (Lindane)	3.66	3.66	3.56	3.76	0.00
Heptachlor	4.01	4.01	3.91	4.11	0.00
Aldrin	4.29	4.29	4.19	4.39	0.00
Heptachlor epoxide	4.79	4.79	4.69	4.89	0.00
Endosulfan I	5.16	5.16	5.06	5.26	0.00
Dieldrin	5.43	5.42	5.32	5.52	-0.01
4,4'-DDE	5.30	5.29	5.19	5.39	-0.01
Endrin	5.70	5.70	5.60	5.80	0.00
Endosulfan II	5.99	5.99	5.89	6.09	0.00
4,4'-DDD	5.85	5.85	5.75	5.95	0.00
Endosulfan sulfate	6.40	6.39	6.29	6.49	0.00
4,4'-DDT	6.10	6.10	6.00	6.20	0.00
Methoxychlor	6.67	6.67	6.57	6.77	0.00
Endrin ketone	6.90	6.90	6.80	7.00	0.00
Endrin aldehyde	6.17	6.17	6.07	6.27	0.00
alpha-Chlordane	5.11	5.10	5.00	5.20	-0.01
gamma-Chlordane	5.04	5.04	4.94	5.14	0.00

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR1 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/28/2025 10/28/2025

Client Sample No.: CCAL06 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PL097715.D **Time Analyzed:** 16:03

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	6.664	6.560	6.760	58.700	50.000	17.4
4,4'-DDE	6.155	6.051	6.251	53.560	50.000	7.1
4,4'-DDT	6.978	6.874	7.074	49.700	50.000	-0.6
Aldrin	5.229	5.127	5.327	51.750	50.000	3.5
alpha-BHC	3.972	3.870	4.070	53.980	50.000	8.0
alpha-Chlordane	5.983	5.881	6.081	52.600	50.000	5.2
beta-BHC	4.488	4.386	4.586	53.670	50.000	7.3
Decachlorobiphenyl	9.000	8.893	9.093	47.600	50.000	-4.8
delta-BHC	4.733	4.631	4.831	54.110	50.000	8.2
Dieldrin	6.303	6.200	6.400	51.850	50.000	3.7
Endosulfan I	6.031	5.928	6.128	52.670	50.000	5.3
Endosulfan II	6.744	6.641	6.841	53.360	50.000	6.7
Endosulfan sulfate	7.107	7.003	7.203	50.710	50.000	1.4
Endrin	6.530	6.427	6.627	50.660	50.000	1.3
Endrin aldehyde	6.873	6.769	6.969	52.540	50.000	5.1
Endrin ketone	7.586	7.481	7.681	51.360	50.000	2.7
gamma-BHC (Lindane)	4.299	4.198	4.398	53.600	50.000	7.2
gamma-Chlordane	5.902	5.801	6.001	51.350	50.000	2.7
Heptachlor	4.891	4.788	4.988	52.360	50.000	4.7
Heptachlor epoxide	5.649	5.547	5.747	52.760	50.000	5.5
Methoxychlor	7.451	7.347	7.547	46.070	50.000	-7.9
Tetrachloro-m-xylene	3.525	3.424	3.624	53.320	50.000	6.6

CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance **Contract:** REMI02
Lab Code: ACE **SDG NO.:** Q3495
GC Column: ZB-MR2 **ID:** 0.32 (mm) **Initi. Calib. Date(s):** 10/28/2025 10/28/2025

Client Sample No.: CCAL06 **Date Analyzed:** 10/31/2025
Lab Sample No.: PSTDCCC050 **Data File :** PL097715.D **Time Analyzed:** 16:03

COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
4,4'-DDD	5.848	5.745	5.945	51.730	50.000	3.5
4,4'-DDE	5.296	5.193	5.393	51.180	50.000	2.4
4,4'-DDT	6.100	5.997	6.197	46.760	50.000	-6.5
Aldrin	4.288	4.187	4.387	52.640	50.000	5.3
alpha-BHC	3.327	3.226	3.426	53.410	50.000	6.8
alpha-Chlordane	5.107	5.004	5.204	50.920	50.000	1.8
beta-BHC	3.955	3.854	4.054	51.150	50.000	2.3
Decachlorobiphenyl	7.984	7.879	8.079	42.150	50.000	-15.7
delta-BHC	4.188	4.086	4.286	52.820	50.000	5.6
Dieldrin	5.426	5.324	5.524	53.500	50.000	7.0
Endosulfan I	5.162	5.059	5.259	49.160	50.000	-1.7
Endosulfan II	5.993	5.890	6.090	49.400	50.000	-1.2
Endosulfan sulfate	6.395	6.291	6.491	48.360	50.000	-3.3
Endrin	5.701	5.599	5.799	52.770	50.000	5.5
Endrin aldehyde	6.171	6.068	6.268	45.730	50.000	-8.5
Endrin ketone	6.899	6.796	6.996	50.160	50.000	0.3
gamma-BHC (Lindane)	3.659	3.558	3.758	53.010	50.000	6.0
gamma-Chlordane	5.043	4.940	5.140	50.840	50.000	1.7
Heptachlor	4.006	3.905	4.105	51.690	50.000	3.4
Heptachlor epoxide	4.791	4.689	4.889	51.480	50.000	3.0
Methoxychlor	6.673	6.569	6.769	45.090	50.000	-9.8
Tetrachloro-m-xylene	2.822	2.721	2.921	51.810	50.000	3.6

PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090648.D Date Analyzed: 10/17/2025

Lab Sample No.(PEM): PEM Time Analyzed: 11:07

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.066	8.970	9.170	22.860	20.000	14.3
Tetrachloro-m-xylene	3.544	3.490	3.590	21.710	20.000	8.6
alpha-BHC	3.994	3.940	4.040	9.050	10.000	-9.5
beta-BHC	4.512	4.460	4.560	10.070	10.000	0.7
gamma-BHC (Lindane)	4.324	4.270	4.370	9.330	10.000	-6.7
Endrin	6.568	6.500	6.640	49.200	50.000	-1.6
4,4'-DDT	7.016	6.950	7.090	100.270	100.000	0.3
Methoxychlor	7.489	7.420	7.560	236.810	250.000	-5.3

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090648.D Date Analyzed: 10/17/2025

Lab Sample No.(PEM): PEM Time Analyzed: 11:07

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.062	7.960	8.160	21.400	20.000	7.0
Tetrachloro-m-xylene	2.875	2.820	2.930	21.220	20.000	6.1
alpha-BHC	3.386	3.340	3.440	10.300	10.000	3.0
beta-BHC	4.019	3.970	4.070	10.460	10.000	4.6
gamma-BHC (Lindane)	3.723	3.670	3.770	10.340	10.000	3.4
Endrin	5.781	5.710	5.850	47.940	50.000	-4.1
4,4'-DDT	6.174	6.100	6.240	95.670	100.000	-4.3
Methoxychlor	6.745	6.670	6.820	201.830	250.000	-19.3

PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090925.D Date Analyzed: 10/31/2025

Lab Sample No.(PEM): PEM Time Analyzed: 10:20

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.070	8.970	9.170	23.600	20.000	18.0
Tetrachloro-m-xylene	3.544	3.490	3.590	23.020	20.000	15.1
alpha-BHC	3.994	3.940	4.040	9.890	10.000	-1.1
beta-BHC	4.513	4.460	4.560	11.020	10.000	10.2
gamma-BHC (Lindane)	4.325	4.270	4.380	10.190	10.000	1.9
Endrin	6.570	6.500	6.640	54.240	50.000	8.5
4,4'-DDT	7.018	6.950	7.090	118.640	100.000	18.6
Methoxychlor	7.491	7.420	7.560	274.190	250.000	9.7

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090925.D Date Analyzed: 10/31/2025

Lab Sample No.(PEM): PEM Time Analyzed: 10:20

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.063	7.960	8.160	30.610	20.000	53.1
Tetrachloro-m-xylene	2.874	2.820	2.920	25.410	20.000	27.1
alpha-BHC	3.386	3.340	3.440	12.590	10.000	25.9
beta-BHC	4.019	3.970	4.070	12.870	10.000	28.7
gamma-BHC (Lindane)	3.723	3.670	3.770	12.620	10.000	26.2
Endrin	5.781	5.710	5.850	59.380	50.000	18.8
4,4'-DDT	6.175	6.100	6.250	132.040	100.000	32.0
Methoxychlor	6.745	6.670	6.820	264.280	250.000	5.7

PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090936.D Date Analyzed: 11/04/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:31

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.062	8.960	9.160	23.370	20.000	16.9
Tetrachloro-m-xylene	3.544	3.490	3.590	21.440	20.000	7.2
alpha-BHC	3.993	3.940	4.040	9.090	10.000	-9.1
beta-BHC	4.511	4.460	4.560	10.360	10.000	3.6
gamma-BHC (Lindane)	4.323	4.270	4.370	9.370	10.000	-6.3
Endrin	6.567	6.500	6.640	51.370	50.000	2.7
4,4'-DDT	7.014	6.940	7.080	113.560	100.000	13.6
Methoxychlor	7.486	7.420	7.560	265.030	250.000	6.0

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 10/17/2025 10/17/2025

Client Sample No. (PEM): PEM - PD090936.D Date Analyzed: 11/04/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:31

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.058	7.960	8.160	28.070	20.000	40.4
Tetrachloro-m-xylene	2.874	2.820	2.920	22.970	20.000	14.9
alpha-BHC	3.385	3.330	3.440	11.220	10.000	12.2
beta-BHC	4.017	3.970	4.070	11.500	10.000	15.0
gamma-BHC (Lindane)	3.721	3.670	3.770	11.340	10.000	13.4
Endrin	5.778	5.710	5.850	52.670	50.000	5.3
4,4'-DDT	6.172	6.100	6.240	117.630	100.000	17.6
Methoxychlor	6.742	6.670	6.810	233.260	250.000	-6.7

PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 10/28/2025 10/28/2025

Client Sample No. (PEM): PEM - PL097637.D Date Analyzed: 10/28/2025

Lab Sample No.(PEM): PEM Time Analyzed: 14:30

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	8.994	8.890	9.090	25.870	20.000	29.4
Tetrachloro-m-xylene	3.524	3.470	3.570	20.090	20.000	0.5
alpha-BHC	3.970	3.920	4.020	8.750	10.000	-12.5
beta-BHC	4.486	4.440	4.540	9.460	10.000	-5.4
gamma-BHC (Lindane)	4.298	4.250	4.350	9.060	10.000	-9.4
Endrin	6.527	6.460	6.600	46.230	50.000	-7.5
4,4'-DDT	6.974	6.900	7.040	94.090	100.000	-5.9
Methoxychlor	7.447	7.380	7.520	224.130	250.000	-10.3

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 10/28/2025 10/28/2025

Client Sample No. (PEM): PEM - PL097637.D Date Analyzed: 10/28/2025

Lab Sample No.(PEM): PEM Time Analyzed: 14:30

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	7.979	7.880	8.080	23.960	20.000	19.8
Tetrachloro-m-xylene	2.822	2.770	2.870	20.030	20.000	0.2
alpha-BHC	3.326	3.280	3.380	8.800	10.000	-12.0
beta-BHC	3.954	3.900	4.000	9.280	10.000	-7.2
gamma-BHC (Lindane)	3.657	3.610	3.710	9.030	10.000	-9.7
Endrin	5.698	5.630	5.770	47.650	50.000	-4.7
4,4'-DDT	6.097	6.030	6.170	97.840	100.000	-2.2
Methoxychlor	6.669	6.600	6.740	208.890	250.000	-16.4

PESTICIDE CALIBRATION VERIFICATION SUMMARY

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

GC Column: ZB-MR1 ID: 0.32 (mm) Initi. Calib. Date(s): 10/28/2025 10/28/2025

Client Sample No. (PEM): PEM - PL097700.D Date Analyzed: 10/31/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:56

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	9.002	8.900	9.100	24.090	20.000	20.5
Tetrachloro-m-xylene	3.524	3.470	3.570	25.470	20.000	27.4
alpha-BHC	3.971	3.920	4.020	11.360	10.000	13.6
beta-BHC	4.488	4.440	4.540	12.130	10.000	21.3
gamma-BHC (Lindane)	4.299	4.250	4.350	11.680	10.000	16.8
Endrin	6.531	6.460	6.600	57.620	50.000	15.2
4,4'-DDT	6.979	6.910	7.050	110.760	100.000	10.8
Methoxychlor	7.452	7.380	7.520	259.100	250.000	3.6

GC Column: ZB-MR2 ID: 0.32 (mm) Initi. Calib. Date(s): 10/28/2025 10/28/2025

Client Sample No. (PEM): PEM - PL097700.D Date Analyzed: 10/31/2025

Lab Sample No.(PEM): PEM Time Analyzed: 09:56

PEM COMPOUND	RT	RT WINDOW		CALC AMOUNT(ng)	NOM AMOUNT(ng)	%D
		FROM	TO			
Decachlorobiphenyl	7.984	7.880	8.080	21.370	20.000	6.9
Tetrachloro-m-xylene	2.822	2.770	2.870	24.530	20.000	22.7
alpha-BHC	3.327	3.280	3.380	11.000	10.000	10.0
beta-BHC	3.956	3.910	4.010	11.540	10.000	15.4
gamma-BHC (Lindane)	3.659	3.610	3.710	11.170	10.000	11.7
Endrin	5.701	5.630	5.770	57.590	50.000	15.2
4,4'-DDT	6.101	6.030	6.170	106.870	100.000	6.9
Methoxychlor	6.674	6.600	6.740	220.400	250.000	-11.8

Analytical Sequence

Client: Remington & Vernick Engineers	SDG No.: Q3495	
Project: Maclearie Park	Instrument ID: ECD_D	
GC Column: ZB-MR1	ID: 0.32 (mm)	Inst. Calib. Date(s): 10/17/2025 10/17/2025

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	10/17/2025	10:53	PD090647.D	9.07	3.55
PEM	PEM	10/17/2025	11:07	PD090648.D	9.07	3.54
RESCHK	RESCHK	10/17/2025	11:20	PD090649.D	9.07	3.54
PSTDICCC100	PSTDICCC100	10/17/2025	11:34	PD090650.D	9.07	3.54
PSTDICCC075	PSTDICCC075	10/17/2025	11:48	PD090651.D	9.07	3.55
PSTDICCC050	PSTDICCC050	10/17/2025	12:01	PD090652.D	9.06	3.55
PSTDICCC025	PSTDICCC025	10/17/2025	12:15	PD090653.D	9.07	3.54
PSTDICCC005	PSTDICCC005	10/17/2025	12:28	PD090654.D	9.07	3.54
PCHLORICC500	PCHLORICC500	10/17/2025	13:09	PD090657.D	9.07	3.55
PTOXICC500	PTOXICC500	10/17/2025	14:17	PD090662.D	9.07	3.55
LBLK	LBLK	10/31/2025	10:06	PD090924.D	9.07	3.54
PEM	PEM	10/31/2025	10:20	PD090925.D	9.07	3.54
PSTDCCC050	PSTDCCC050	10/31/2025	11:09	PD090926.D	9.08	3.55
PB170348BL	PB170348BL	10/31/2025	14:37	PD090927.D	9.08	3.55
TW-1025-1	Q3495-03	10/31/2025	15:34	PD090931.D	9.07	3.54
LBLK	LBLK	10/31/2025	15:48	PD090932.D	9.07	3.55
PSTDCCC050	PSTDCCC050	10/31/2025	16:01	PD090933.D	9.07	3.55
PEM	PEM	11/04/2025	09:31	PD090936.D	9.06	3.54
LBLK	LBLK	11/04/2025	12:35	PD090948.D	9.06	3.54
PSTDCCC050	PSTDCCC050	11/04/2025	13:29	PD090949.D	9.07	3.55
PB170348BSD	PB170348BSD	11/04/2025	13:52	PD090950.D	9.07	3.55
PB170348BS	PB170348BS	11/04/2025	14:22	PD090952.D	9.07	3.54
LBLK	LBLK	11/04/2025	16:31	PD090956.D	9.06	3.55
PSTDCCC050	PSTDCCC050	11/04/2025	16:44	PD090957.D	9.06	3.54
LBLK	LBLK	10/28/2025	14:17	PL097636.D	8.99	3.52
PEM	PEM	10/28/2025	14:30	PL097637.D	8.99	3.52
RESCHK	RESCHK	10/28/2025	14:44	PL097638.D	8.99	3.52
PSTDICCC100	PSTDICCC100	10/28/2025	14:57	PL097639.D	8.99	3.52
PSTDICCC075	PSTDICCC075	10/28/2025	15:11	PL097640.D	8.99	3.52
PSTDICCC050	PSTDICCC050	10/28/2025	15:25	PL097641.D	8.99	3.52
PSTDICCC025	PSTDICCC025	10/28/2025	15:38	PL097642.D	9.00	3.52
PSTDICCC005	PSTDICCC005	10/28/2025	15:52	PL097643.D	8.99	3.52
PCHLORICC500	PCHLORICC500	10/28/2025	16:32	PL097646.D	8.99	3.52
PTOXICC500	PTOXICC500	10/28/2025	17:40	PL097651.D	8.99	3.52
LBLK	LBLK	10/31/2025	09:42	PL097699.D	9.00	3.52
PEM	PEM	10/31/2025	09:56	PL097700.D	9.00	3.52
PSTDCCC050	PSTDCCC050	10/31/2025	11:02	PL097701.D	9.01	3.52
PB170340BL	PB170340BL	10/31/2025	13:06	PL097704.D	9.01	3.52
PB170340BS	PB170340BS	10/31/2025	13:19	PL097705.D	9.00	3.52
SB-1025-2	Q3495-02	10/31/2025	13:33	PL097706.D	9.00	3.52
SB-1025-2MS	Q3495-02MS	10/31/2025	13:47	PL097707.D	9.00	3.52
SB-1025-2MSD	Q3495-02MSD	10/31/2025	14:00	PL097708.D	9.00	3.52

Analytical Sequence

I.BLK	I.BLK	10/31/2025	15:49	PL097714.D	9.00	3.53
PSTDCCC050	PSTDCCC050	10/31/2025	16:03	PL097715.D	9.00	3.53

Analytical Sequence

Client: Remington & Vernick Engineers	SDG No.: Q3495	
Project: Maclearie Park	Instrument ID: ECD_D	
GC Column: ZB-MR2	ID: 0.32 (mm)	Inst. Calib. Date(s): 10/17/2025 10/17/2025

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	10/17/2025	10:53	PD090647.D	8.06	2.88
PEM	PEM	10/17/2025	11:07	PD090648.D	8.06	2.88
RESCHK	RESCHK	10/17/2025	11:20	PD090649.D	8.06	2.87
PSTDICCC100	PSTDICCC100	10/17/2025	11:34	PD090650.D	8.06	2.87
PSTDICCC075	PSTDICCC075	10/17/2025	11:48	PD090651.D	8.06	2.88
PSTDICCC050	PSTDICCC050	10/17/2025	12:01	PD090652.D	8.06	2.88
PSTDICCC025	PSTDICCC025	10/17/2025	12:15	PD090653.D	8.06	2.87
PSTDICCC005	PSTDICCC005	10/17/2025	12:28	PD090654.D	8.06	2.87
PCHLORICC500	PCHLORICC500	10/17/2025	13:09	PD090657.D	8.06	2.87
PTOXICC500	PTOXICC500	10/17/2025	14:17	PD090662.D	8.06	2.88
IBLK	IBLK	10/31/2025	10:06	PD090924.D	8.06	2.87
PEM	PEM	10/31/2025	10:20	PD090925.D	8.06	2.87
PSTDCCC050	PSTDCCC050	10/31/2025	11:09	PD090926.D	8.07	2.87
PB170348BL	PB170348BL	10/31/2025	14:37	PD090927.D	8.07	2.87
TW-1025-1	Q3495-03	10/31/2025	15:34	PD090931.D	8.06	2.87
IBLK	IBLK	10/31/2025	15:48	PD090932.D	8.06	2.88
PSTDCCC050	PSTDCCC050	10/31/2025	16:01	PD090933.D	8.06	2.87
PEM	PEM	11/04/2025	09:31	PD090936.D	8.06	2.87
IBLK	IBLK	11/04/2025	12:35	PD090948.D	8.06	2.87
PSTDCCC050	PSTDCCC050	11/04/2025	13:29	PD090949.D	8.06	2.87
PB170348BSD	PB170348BSD	11/04/2025	13:52	PD090950.D	8.06	2.87
PB170348BS	PB170348BS	11/04/2025	14:22	PD090952.D	8.06	2.87
IBLK	IBLK	11/04/2025	16:31	PD090956.D	8.06	2.87
PSTDCCC050	PSTDCCC050	11/04/2025	16:44	PD090957.D	8.06	2.87
IBLK	IBLK	10/28/2025	14:17	PL097636.D	7.98	2.82
PEM	PEM	10/28/2025	14:30	PL097637.D	7.98	2.82
RESCHK	RESCHK	10/28/2025	14:44	PL097638.D	7.98	2.82
PSTDICCC100	PSTDICCC100	10/28/2025	14:57	PL097639.D	7.98	2.82
PSTDICCC075	PSTDICCC075	10/28/2025	15:11	PL097640.D	7.98	2.82
PSTDICCC050	PSTDICCC050	10/28/2025	15:25	PL097641.D	7.98	2.82
PSTDICCC025	PSTDICCC025	10/28/2025	15:38	PL097642.D	7.98	2.82
PSTDICCC005	PSTDICCC005	10/28/2025	15:52	PL097643.D	7.98	2.82
PCHLORICC500	PCHLORICC500	10/28/2025	16:32	PL097646.D	7.98	2.82
PTOXICC500	PTOXICC500	10/28/2025	17:40	PL097651.D	7.98	2.82
IBLK	IBLK	10/31/2025	09:42	PL097699.D	7.98	2.82
PEM	PEM	10/31/2025	09:56	PL097700.D	7.98	2.82
PSTDCCC050	PSTDCCC050	10/31/2025	11:02	PL097701.D	7.98	2.82
PB170340BL	PB170340BL	10/31/2025	13:06	PL097704.D	7.99	2.82
PB170340BS	PB170340BS	10/31/2025	13:19	PL097705.D	7.98	2.82
SB-1025-2	Q3495-02	10/31/2025	13:33	PL097706.D	7.98	2.82
SB-1025-2MS	Q3495-02MS	10/31/2025	13:47	PL097707.D	7.98	2.82
SB-1025-2MSD	Q3495-02MSD	10/31/2025	14:00	PL097708.D	7.98	2.82

Analytical Sequence

I.BLK	I.BLK	10/31/2025	15:49	PL097714.D	7.98	2.82
PSTDCCC050	PSTDCCC050	10/31/2025	16:03	PL097715.D	7.98	2.82

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170340BS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170340BS

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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	6.75	6.70	6.80	17.0	9.9
	2	5.99	5.94	6.04	15.4	
4,4'-DDD	1	6.67	6.62	6.72	18.0	11.1
	2	5.85	5.80	5.90	16.1	
4,4'-DDT	1	6.98	6.93	7.03	16.3	9
	2	6.10	6.05	6.15	14.9	
Endrin aldehyde	1	6.87	6.82	6.92	16.3	13.8
	2	6.17	6.12	6.22	14.2	
Endosulfan sulfate	1	7.11	7.06	7.16	16.4	8.9
	2	6.39	6.34	6.44	15.0	
Methoxychlor	1	7.45	7.40	7.50	15.2	6.8
	2	6.67	6.62	6.72	14.2	
Endrin ketone	1	7.59	7.54	7.64	16.8	4.3
	2	6.90	6.85	6.95	16.1	
alpha-BHC	1	3.97	3.92	4.02	17.0	0.6
	2	3.33	3.28	3.38	16.9	
gamma-BHC (Lindane)	1	4.30	4.25	4.35	17.0	0.6
	2	3.66	3.61	3.71	16.9	
Heptachlor	1	4.89	4.84	4.94	17.1	4.2
	2	4.01	3.96	4.06	16.4	
Aldrin	1	5.23	5.18	5.28	16.6	1.2
	2	4.29	4.24	4.34	16.8	
beta-BHC	1	4.49	4.44	4.54	17.1	7.3
	2	3.95	3.90	4.00	15.9	
delta-BHC	1	4.73	4.68	4.78	17.4	4.1
	2	4.19	4.14	4.24	16.7	
Heptachlor epoxide	1	5.65	5.60	5.70	16.0	1.9
	2	4.79	4.74	4.84	16.3	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170340BS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170340BS

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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	6.03	5.98	6.08	16.1	1.9
	2	5.16	5.11	5.21	15.8	
gamma-Chlordane	1	5.90	5.85	5.95	16.5	3.1
	2	5.04	4.99	5.09	16.0	
alpha-Chlordane	1	5.98	5.93	6.03	16.2	0.6
	2	5.11	5.06	5.16	16.1	
4,4'-DDE	1	6.15	6.10	6.20	16.8	3
	2	5.30	5.25	5.35	16.3	
Dieldrin	1	6.30	6.25	6.35	16.1	2.5
	2	5.43	5.38	5.48	16.5	
Endrin	1	6.53	6.48	6.58	16.0	7.8
	2	5.70	5.65	5.75	17.3	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170348BS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170348BS

Date(s) Analyzed: 11/04/2025

11/04/2025

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	0.50	1.2
	2	5.92	5.87	5.97	0.49	
4,4'-DDE	1	6.19	6.14	6.24	0.48	3.7
	2	5.36	5.31	5.41	0.50	
4,4'-DDT	1	7.02	6.97	7.07	0.48	2
	2	6.17	6.12	6.22	0.47	
alpha-BHC	1	3.99	3.94	4.04	0.52	1.6
	2	3.39	3.34	3.44	0.51	
Aldrin	1	5.26	5.21	5.31	0.51	1.5
	2	4.36	4.31	4.41	0.51	
alpha-Chlordane	1	6.02	5.97	6.07	0.49	3.1
	2	5.18	5.13	5.23	0.50	
Endosulfan II	1	6.78	6.73	6.83	0.48	2.8
	2	6.07	6.02	6.12	0.47	
Endosulfan sulfate	1	7.14	7.09	7.19	0.46	0.6
	2	6.47	6.42	6.52	0.46	
beta-BHC	1	4.51	4.46	4.56	0.50	0.9
	2	4.02	3.97	4.07	0.50	
delta-BHC	1	4.76	4.71	4.81	0.51	1.1
	2	4.25	4.20	4.30	0.51	
Endosulfan I	1	6.07	6.02	6.12	0.49	2.9
	2	5.24	5.19	5.29	0.50	
Dieldrin	1	6.34	6.29	6.39	0.49	1
	2	5.50	5.45	5.55	0.49	
Endrin aldehyde	1	6.91	6.86	6.96	0.45	0.4
	2	6.25	6.20	6.30	0.45	
Methoxychlor	1	7.49	7.44	7.54	0.48	6.2
	2	6.74	6.69	6.79	0.45	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170348BS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170348BS

Date(s) Analyzed: 11/04/2025 11/04/2025

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 ketone	1	7.63	7.58	7.68	0.46	4.9
	2	6.98	6.93	7.03	0.44	
gamma-BHC (Lindane)	1	4.32	4.27	4.37	0.51	0.6
	2	3.72	3.67	3.77	0.51	
Heptachlor	1	4.92	4.87	4.97	0.63	16.3
	2	4.07	4.02	4.12	0.54	
Heptachlor epoxide	1	5.68	5.63	5.73	0.50	0.7
	2	4.86	4.81	4.91	0.50	
gamma-Chlordane	1	5.94	5.89	5.99	0.50	0.6
	2	5.12	5.07	5.17	0.50	
Endrin	1	6.57	6.52	6.62	0.48	1.3
	2	5.78	5.73	5.83	0.48	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170348BSD

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170348BSD

Date(s) Analyzed: 11/04/2025 11/04/2025

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	0.49	3.8
	2	5.92	5.87	5.97	0.47	
4,4'-DDT	1	7.02	6.97	7.07	0.48	3.2
	2	6.17	6.12	6.22	0.46	
alpha-BHC	1	4.00	3.95	4.05	0.52	1.9
	2	3.39	3.34	3.44	0.51	
Aldrin	1	5.27	5.22	5.32	0.51	1.3
	2	4.36	4.31	4.41	0.50	
beta-BHC	1	4.52	4.47	4.57	0.49	2
	2	4.02	3.97	4.07	0.50	
alpha-Chlordane	1	6.03	5.98	6.08	0.47	4.7
	2	5.18	5.13	5.23	0.49	
4,4'-DDE	1	6.20	6.15	6.25	0.48	0.4
	2	5.37	5.32	5.42	0.48	
Endosulfan II	1	6.79	6.74	6.84	0.48	5.2
	2	6.07	6.02	6.12	0.45	
Endrin aldehyde	1	6.92	6.87	6.97	0.44	0.2
	2	6.25	6.20	6.30	0.45	
Endosulfan sulfate	1	7.15	7.10	7.20	0.45	3.3
	2	6.47	6.42	6.52	0.44	
Methoxychlor	1	7.50	7.45	7.55	0.47	7.6
	2	6.74	6.69	6.79	0.43	
Endrin ketone	1	7.63	7.58	7.68	0.46	10.3
	2	6.98	6.93	7.03	0.41	
gamma-BHC (Lindane)	1	4.33	4.28	4.38	0.51	1.5
	2	3.72	3.67	3.77	0.51	
Heptachlor	1	4.93	4.88	4.98	0.63	17.5
	2	4.07	4.02	4.12	0.53	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

PB170348BSD

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: PB170348BSD

Date(s) Analyzed: 11/04/2025 11/04/2025

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	0.51	1.6
	2	4.25	4.20	4.30	0.50	
Heptachlor epoxide	1	5.69	5.64	5.74	0.49	2.3
	2	4.86	4.81	4.91	0.50	
Endosulfan I	1	6.08	6.03	6.13	0.48	1.3
	2	5.24	5.19	5.29	0.49	
gamma-Chlordane	1	5.95	5.90	6.00	0.48	1.9
	2	5.12	5.07	5.17	0.49	
Dieldrin	1	6.35	6.30	6.40	0.49	1.5
	2	5.50	5.45	5.55	0.48	
Endrin	1	6.58	6.53	6.63	0.47	3.4
	2	5.78	5.73	5.83	0.46	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

SB-1025-2MS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: Q3495-02MS

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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	6.74	6.69	6.79	21.9	4.2
	2	5.99	5.94	6.04	21.0	
4,4'-DDD	1	6.66	6.61	6.71	23.3	13.8
	2	5.85	5.80	5.90	20.3	
4,4'-DDT	1	6.98	6.93	7.03	22.1	7
	2	6.10	6.05	6.15	20.6	
Endrin aldehyde	1	6.87	6.82	6.92	21.6	11.8
	2	6.17	6.12	6.22	19.2	
Endosulfan sulfate	1	7.11	7.06	7.16	21.9	8.1
	2	6.40	6.35	6.45	20.2	
Methoxychlor	1	7.45	7.40	7.50	21.5	9.2
	2	6.67	6.62	6.72	19.6	
Endrin ketone	1	7.59	7.54	7.64	21.7	8.7
	2	6.90	6.85	6.95	19.9	
alpha-BHC	1	3.97	3.92	4.02	22.1	1.8
	2	3.33	3.28	3.38	21.7	
gamma-BHC (Lindane)	1	4.30	4.25	4.35	22.1	1.4
	2	3.66	3.61	3.71	21.8	
Heptachlor	1	4.89	4.84	4.94	22.9	6.8
	2	4.01	3.96	4.06	21.4	
Aldrin	1	5.23	5.18	5.28	21.7	0.5
	2	4.29	4.24	4.34	21.8	
beta-BHC	1	4.49	4.44	4.54	22.5	8.3
	2	3.96	3.91	4.01	20.7	
delta-BHC	1	4.73	4.68	4.78	23.0	5.8
	2	4.19	4.14	4.24	21.7	
Heptachlor epoxide	1	5.65	5.60	5.70	21.1	1.9
	2	4.79	4.74	4.84	21.5	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

SB-1025-2MS

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: Q3495-02MS

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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	6.03	5.98	6.08	21.4	0.9
	2	5.16	5.11	5.21	21.2	
gamma-Chlordane	1	5.90	5.85	5.95	21.9	3.7
	2	5.04	4.99	5.09	21.1	
alpha-Chlordane	1	5.98	5.93	6.03	21.4	0.5
	2	5.11	5.06	5.16	21.3	
4,4'-DDE	1	6.16	6.11	6.21	22.6	5
	2	5.30	5.25	5.35	21.5	
Dieldrin	1	6.30	6.25	6.35	21.5	0.5
	2	5.43	5.38	5.48	21.6	
Endrin	1	6.53	6.48	6.58	21.2	4.2
	2	5.70	5.65	5.75	22.1	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

SB-1025-2MSD

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: Q3495-02MSD

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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 sulfate	1	7.11	7.06	7.16	22.3	7
	2	6.40	6.35	6.45	20.8	
Methoxychlor	1	7.45	7.40	7.50	21.4	6.8
	2	6.67	6.62	6.72	20.0	
Endrin ketone	1	7.59	7.54	7.64	22.3	8.4
	2	6.90	6.85	6.95	20.5	
alpha-BHC	1	3.97	3.92	4.02	22.7	1.8
	2	3.33	3.28	3.38	22.3	
gamma-BHC (Lindane)	1	4.30	4.25	4.35	22.6	0.9
	2	3.66	3.61	3.71	22.4	
Heptachlor	1	4.89	4.84	4.94	23.5	8
	2	4.01	3.96	4.06	21.7	
Aldrin	1	5.23	5.18	5.28	22.2	0.9
	2	4.29	4.24	4.34	22.4	
beta-BHC	1	4.49	4.44	4.54	23.0	8.1
	2	3.96	3.91	4.01	21.2	
delta-BHC	1	4.73	4.68	4.78	23.6	5.7
	2	4.19	4.14	4.24	22.3	
Heptachlor epoxide	1	5.65	5.60	5.70	21.6	1.4
	2	4.79	4.74	4.84	21.9	
Endosulfan I	1	6.03	5.98	6.08	22.1	1.8
	2	5.16	5.11	5.21	21.7	
gamma-Chlordane	1	5.90	5.85	5.95	22.3	3.7
	2	5.04	4.99	5.09	21.5	
alpha-Chlordane	1	5.98	5.93	6.03	22.1	1.8
	2	5.11	5.06	5.16	21.7	
4,4'-DDE	1	6.16	6.11	6.21	23.3	5.7
	2	5.30	5.25	5.35	22.0	

COMPOUND DETECTION SUMMARY

CLIENT SAMPLE NO.

SB-1025-2MSD

Lab Name: Alliance

Contract: REMI02

Lab Code: ACE

SDG NO.: Q3495

Lab Sample ID: Q3495-02MSD

Date(s) Analyzed: 10/31/2025 10/31/2025

Instrument ID (1): ECD_L

Instrument ID (2): ECD_L

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		
Dieldrin	1	6.30	6.25	6.35	22.0	0
	2	5.43	5.38	5.48	22.0	
Endrin	1	6.53	6.48	6.58	21.7	3.6
	2	5.70	5.65	5.75	22.5	
Endosulfan II	1	6.75	6.70	6.80	22.4	4.1
	2	5.99	5.94	6.04	21.5	
4,4'-DDD	1	6.67	6.62	6.72	23.8	13
	2	5.85	5.80	5.90	20.9	
4,4'-DDT	1	6.98	6.93	7.03	22.4	6.5
	2	6.10	6.05	6.15	21.0	
Endrin aldehyde	1	6.87	6.82	6.92	22.2	11.9
	2	6.17	6.12	6.22	19.7	



SAMPLE RAW DATA

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
 Data File : PL097706.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 13:33
 Operator : AR\AJ
 Sample : Q3495-02
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 ECD_L
 ClientSampleId :
 SB-1025-2

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 11/03/2025
 Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:54:10 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
 Quant Title : GC Extractables
 QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

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

System Monitoring Compounds						
1) SA Tetrachlo...	3.522	2.821	80750731	113.5E6	19.734m	18.859
28) SA Decachlor...	9.002	7.984	47978033	76489819	16.126	14.311

Target Compounds

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

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
Data File : PL097706.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 13:33
Operator : AR\AJ
Sample : Q3495-02
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :

ECD_L

ClientSampleId :

SB-1025-2

Manual Integrations

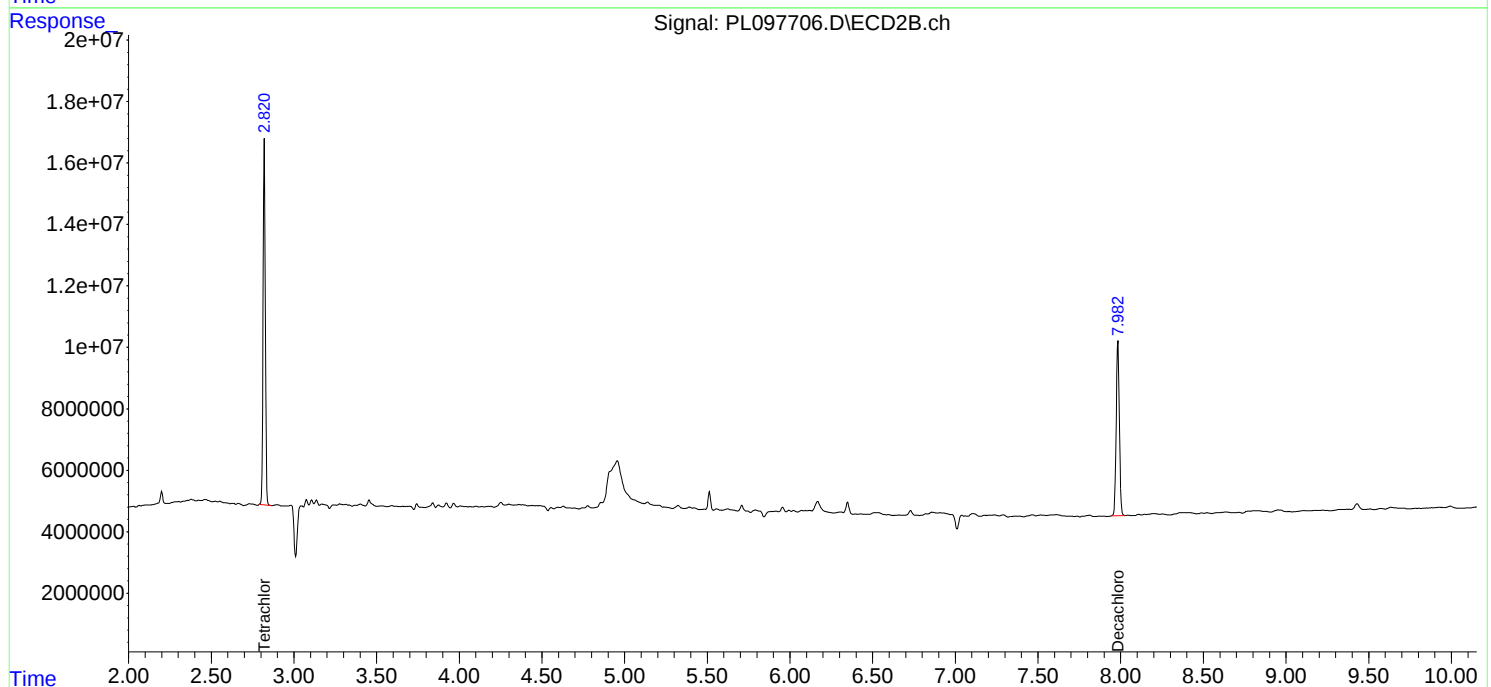
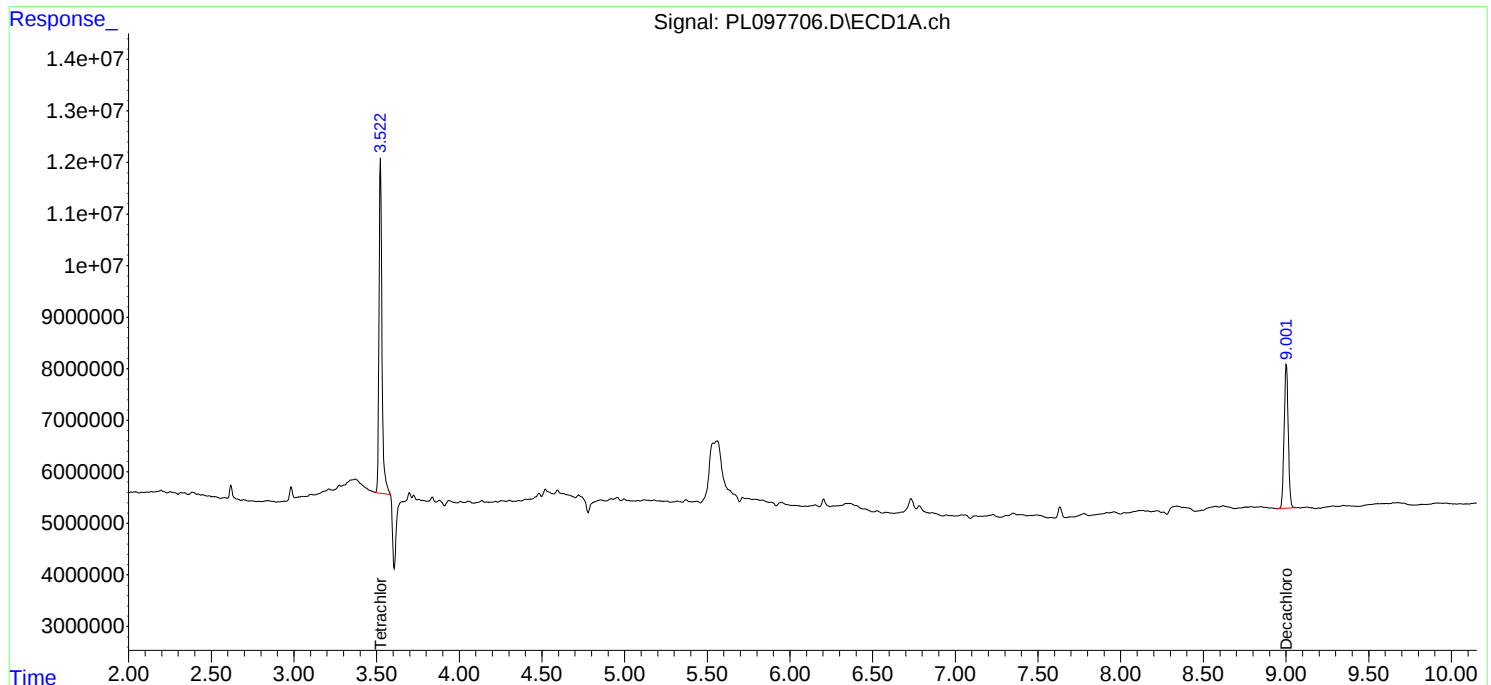
APPROVED

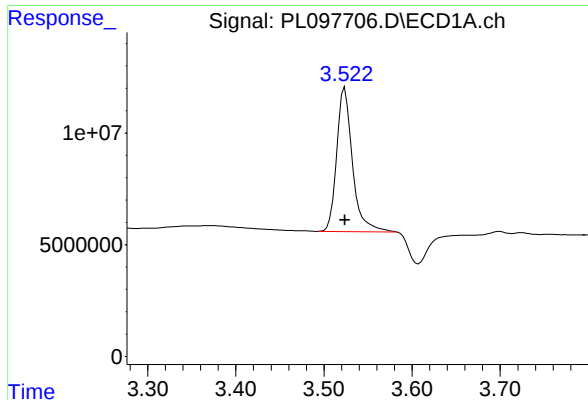
Reviewed By :Rahul Chavli 11/03/2025

Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:54:10 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
Quant Title : GC Extractables
QLast Update : Tue Oct 28 17:14:31 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 x0.25µm





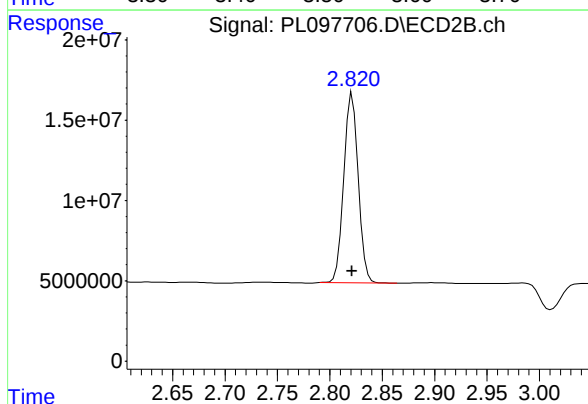
#1 Tetrachloro-m-xylene

R.T.: 3.522 min
Delta R.T.: -0.002 min
Response: 80750731
Conc: 19.73 ng/ml

Instrument :
ECD_L
ClientSampleId :
SB-1025-2

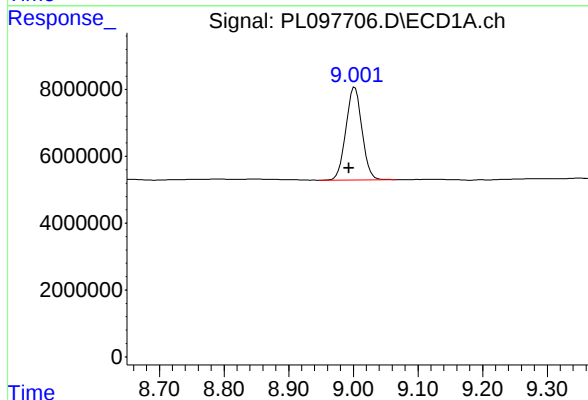
Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 11/03/2025
Supervised By :Yogesh Patel 11/03/2025



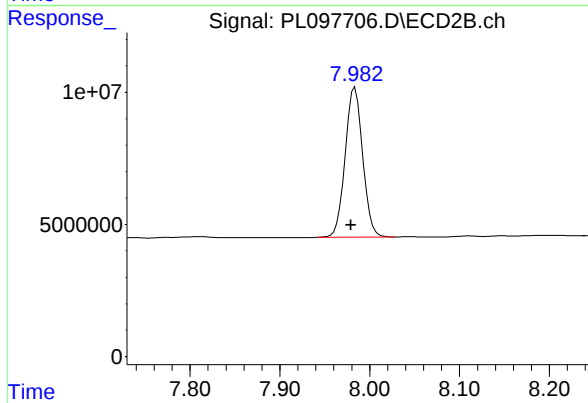
#1 Tetrachloro-m-xylene

R.T.: 2.821 min
Delta R.T.: 0.000 min
Response: 113511316
Conc: 18.86 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.002 min
Delta R.T.: 0.009 min
Response: 47978033
Conc: 16.13 ng/ml



#28 Decachlorobiphenyl

R.T.: 7.984 min
Delta R.T.: 0.005 min
Response: 76489819
Conc: 14.31 ng/ml

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD103125\
 Data File : PD090931.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 15:34
 Operator : AR\AJ
 Sample : Q3495-03
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 TW-1025-1

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 11/03/2025
 Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:54:04 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
 Quant Title : GC Extractables
 QLast Update : Fri Oct 17 17:43:28 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 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...	3.543	2.874	51572972	592.4E6	16.561m	19.896
28) SA Decachlor...	9.070	8.061	62821856	509.8E6	13.429	16.829 #

Target Compounds

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

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD103125\
Data File : PD090931.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 15:34
Operator : AR\AJ
Sample : Q3495-03
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

TW-1025-1

Manual Integrations

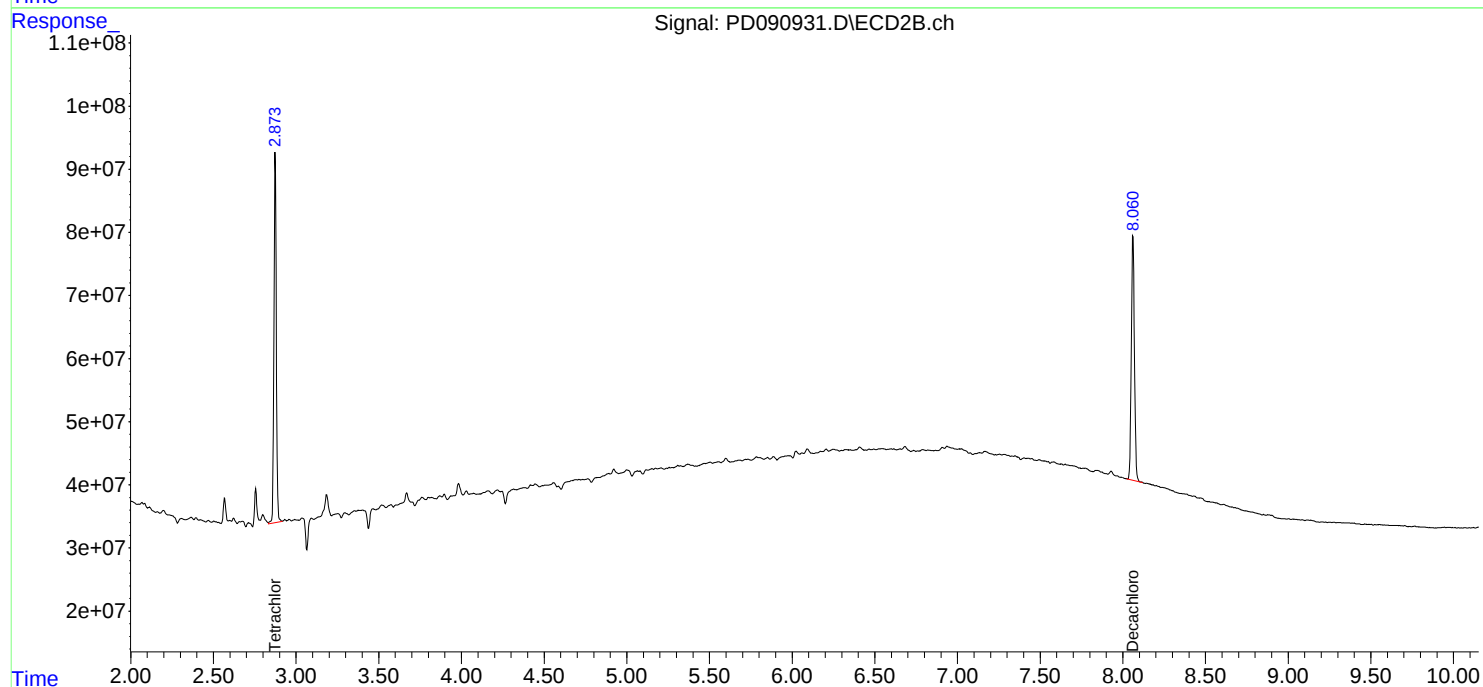
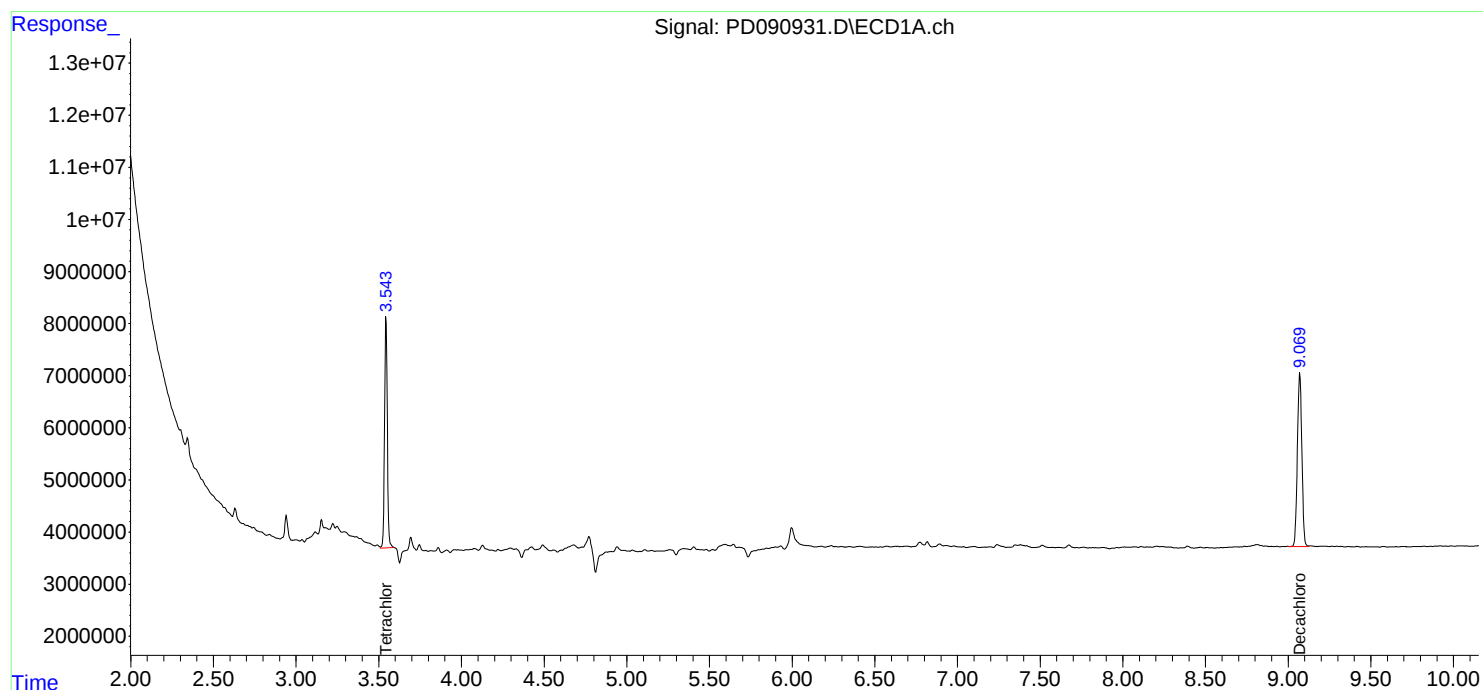
APPROVED

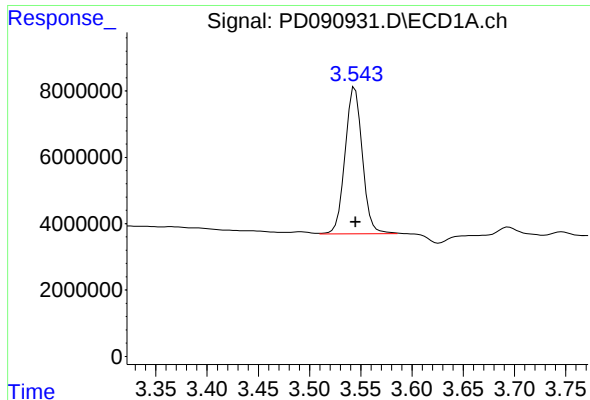
Reviewed By :Rahul Chavli 11/03/2025

Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:54:04 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
Quant Title : GC Extractables
QLast Update : Fri Oct 17 17:43:28 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 x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm





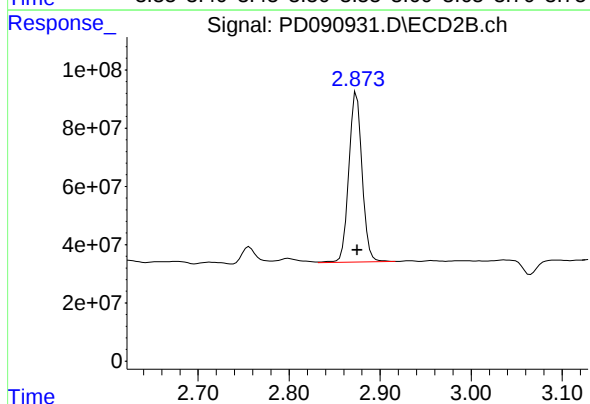
#1 Tetrachloro-m-xylene

R.T.: 3.543 min
Delta R.T.: -0.002 min
Response: 51572972
Conc: 16.56 ng/ml

Instrument :
ECD_D
ClientSampleId :
TW-1025-1

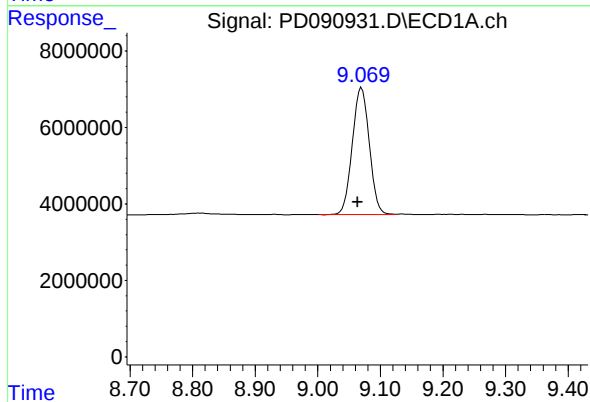
Manual Integrations
APPROVED

Reviewed By :Rahul Chavli 11/03/2025
Supervised By :Yogesh Patel 11/03/2025



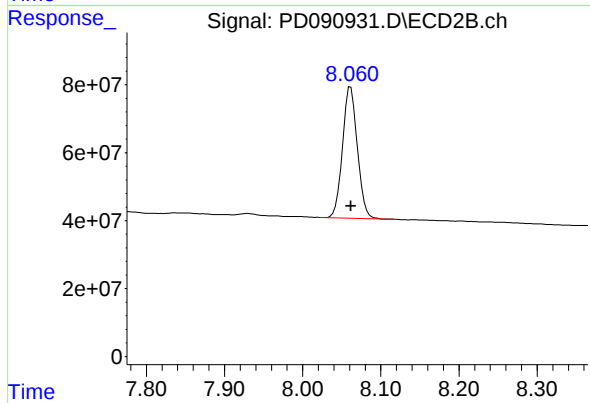
#1 Tetrachloro-m-xylene

R.T.: 2.874 min
Delta R.T.: 0.000 min
Response: 592433104
Conc: 19.90 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.070 min
Delta R.T.: 0.006 min
Response: 62821856
Conc: 13.43 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.061 min
Delta R.T.: 0.000 min
Response: 509774537
Conc: 16.83 ng/ml

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
 Data File : PL097704.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 13:06
 Operator : AR\AJ
 Sample : PB170340BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 ECD_L
 ClientSampleId :
 PB170340BL

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:53:37 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
 Quant Title : GC Extractables
 QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

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

System Monitoring Compounds						
1) SA Tetrachlo...	3.524	2.818	79274682	120.3E6	19.374	19.985
28) SA Decachlor...	9.007	7.985	56171213	85245821	18.880	15.949

Target Compounds

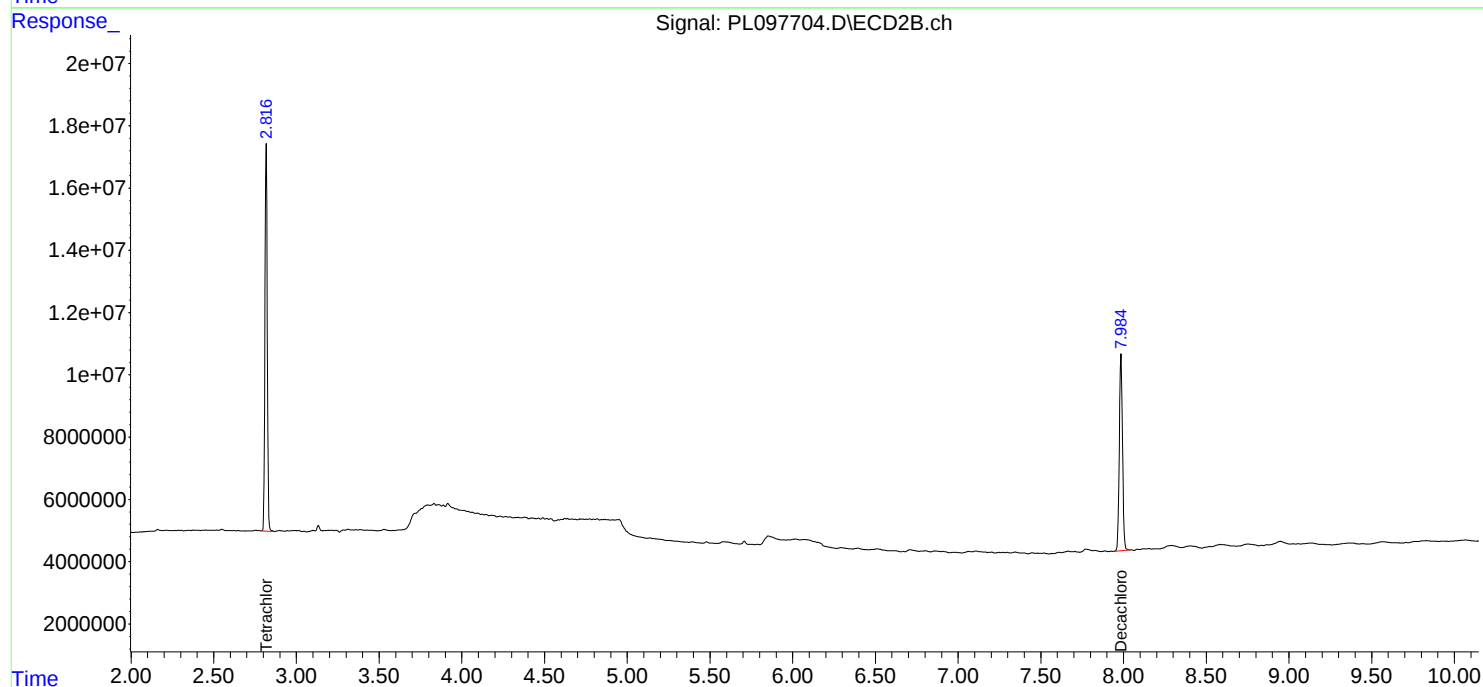
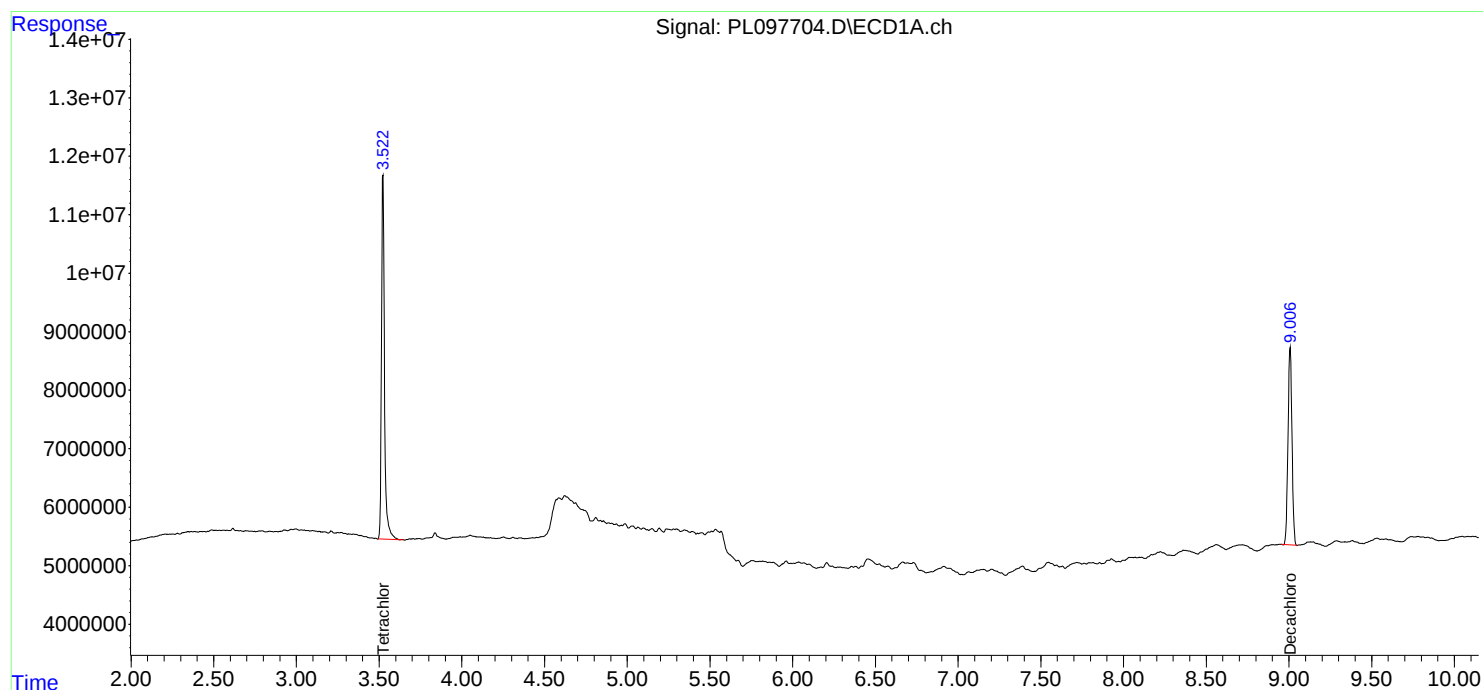
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

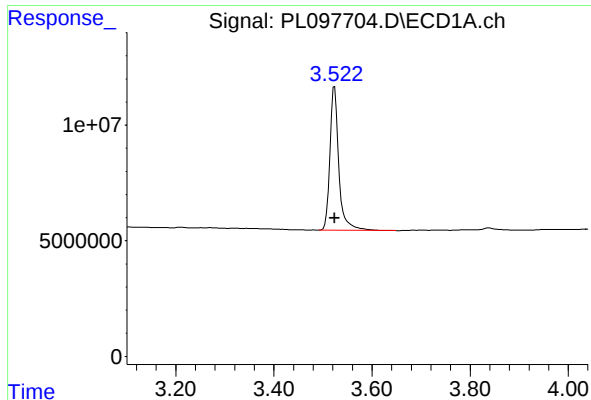
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
Data File : PL097704.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 13:06
Operator : AR\AJ
Sample : PB170340BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
ECD_L
ClientSampleId :
PB170340BL

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:53:37 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
Quant Title : GC Extractables
QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

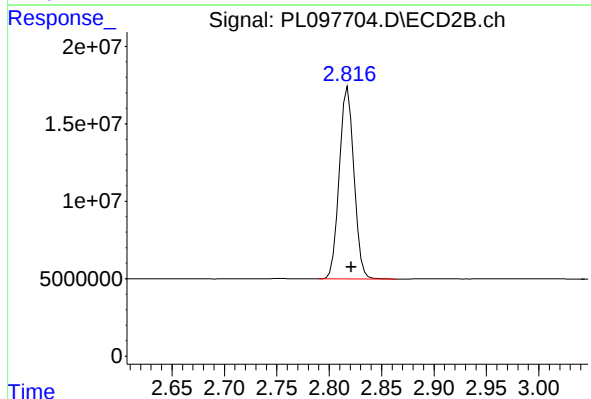




#1 Tetrachloro-m-xylene

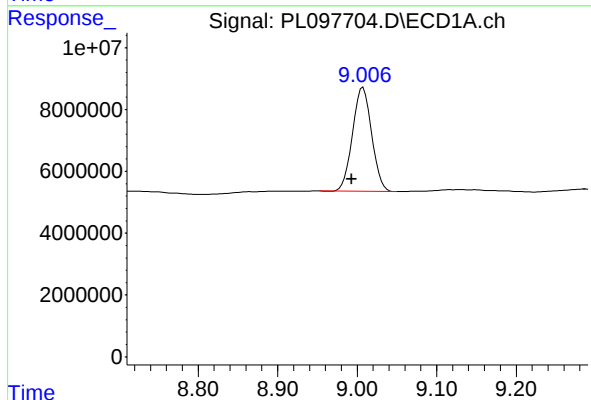
R.T.: 3.524 min
Delta R.T.: 0.000 min
Response: 79274682
Conc: 19.37 ng/ml

Instrument :
ECD_L
ClientSampleId :
PB170340BL



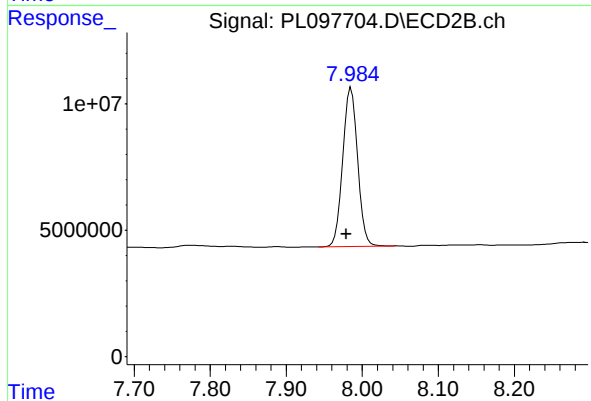
#1 Tetrachloro-m-xylene

R.T.: 2.818 min
Delta R.T.: -0.003 min
Response: 120288529
Conc: 19.98 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.007 min
Delta R.T.: 0.014 min
Response: 56171213
Conc: 18.88 ng/ml



#28 Decachlorobiphenyl

R.T.: 7.985 min
Delta R.T.: 0.006 min
Response: 85245821
Conc: 15.95 ng/ml

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD103125\
 Data File : PD090927.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 14:37
 Operator : AR\AJ
 Sample : PB170348BL
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB170348BL

A
B
C
D
E
F
G
H
I
J
K
L

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:52:42 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
 Quant Title : GC Extractables
 QLast Update : Fri Oct 17 17:43:28 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 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...	3.552	2.873	58389156	667.3E6	18.750	22.409
28) SA Decachlor...	9.083	8.066	102.7E6	823.9E6	21.950	27.200

Target Compounds

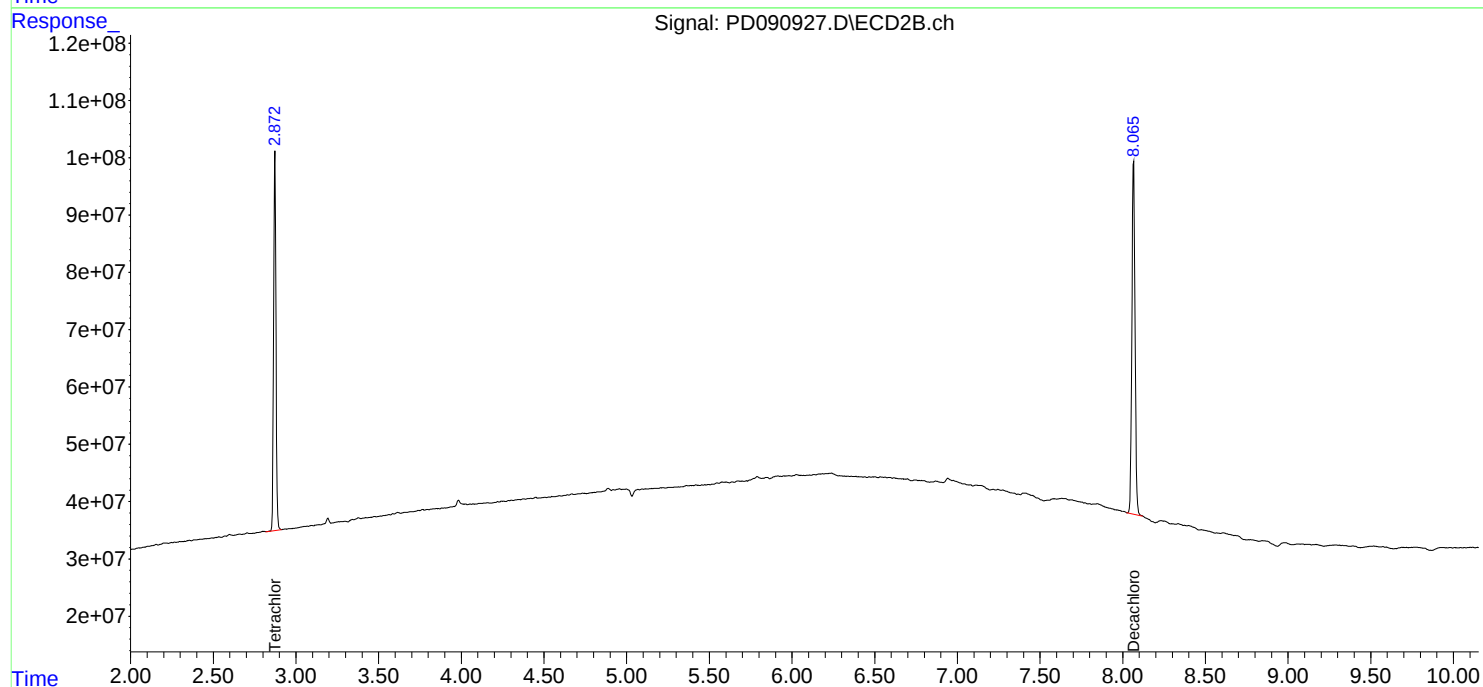
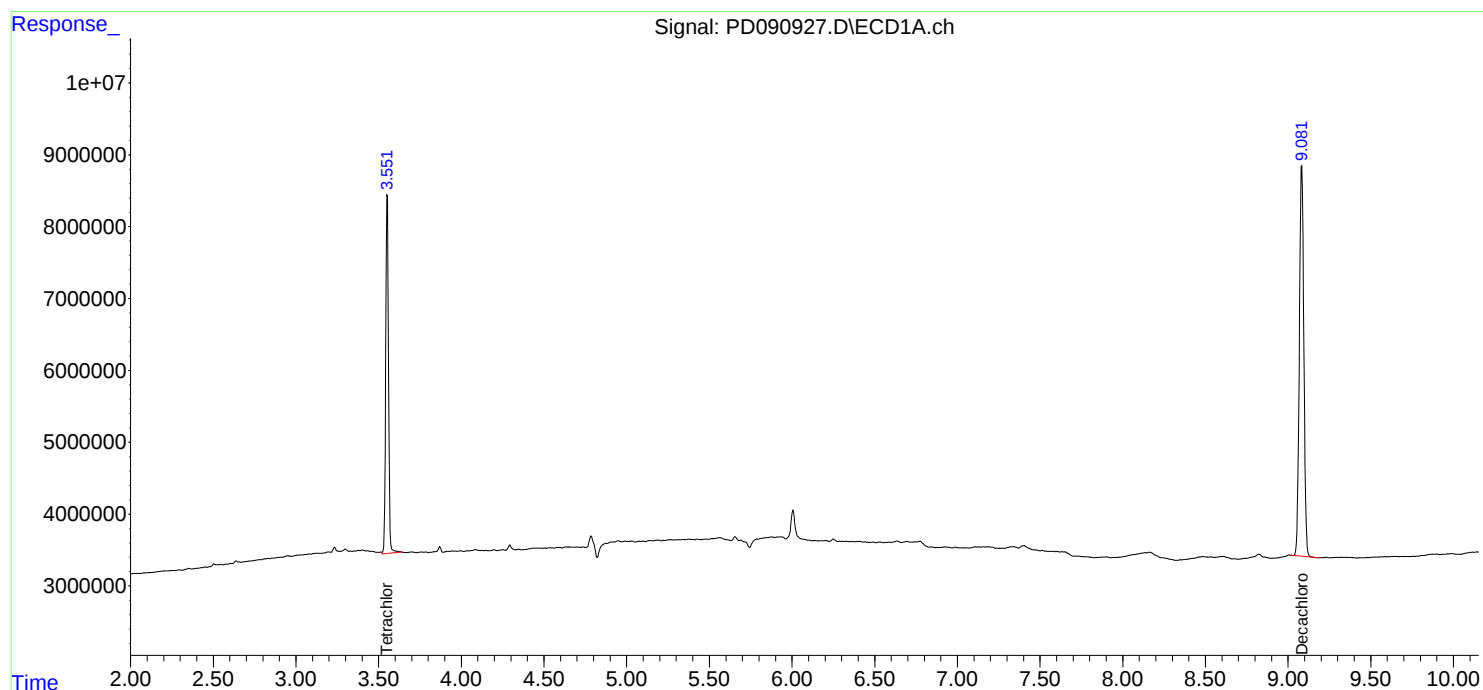
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

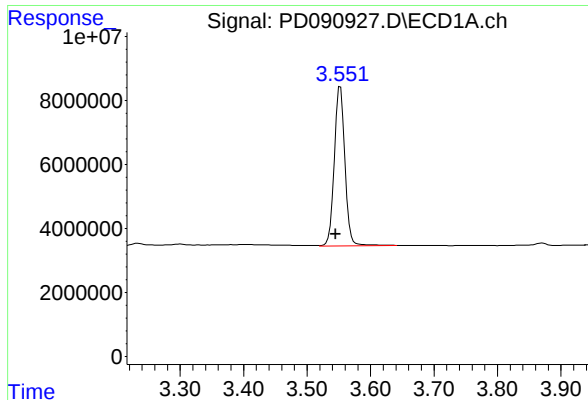
Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD103125\
Data File : PD090927.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 14:37
Operator : AR\AJ
Sample : PB170348BL
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PB170348BL

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:52:42 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
Quant Title : GC Extractables
QLast Update : Fri Oct 17 17:43:28 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 0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm

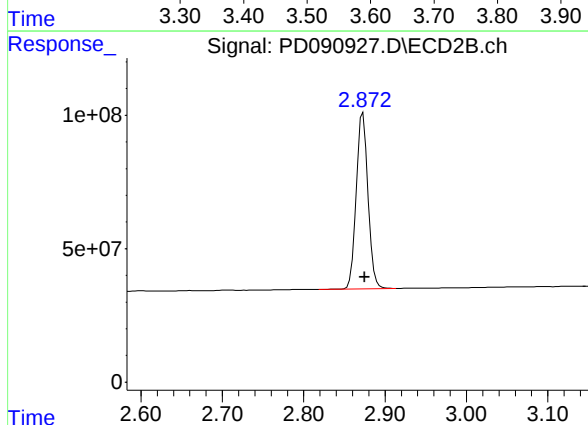




#1 Tetrachloro-m-xylene

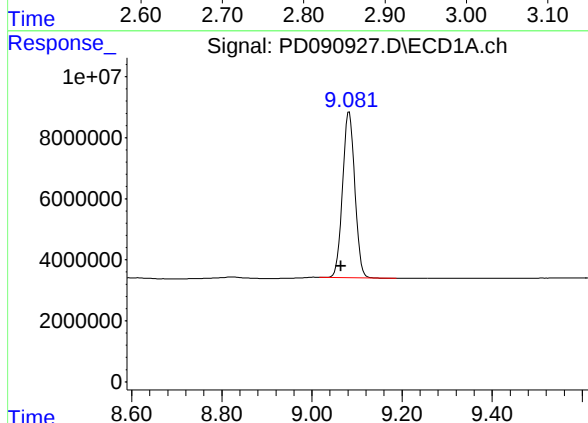
R.T.: 3.552 min
Delta R.T.: 0.007 min
Response: 58389156
Conc: 18.75 ng/ml

Instrument :
ECD_D
ClientSampleId :
PB170348BL



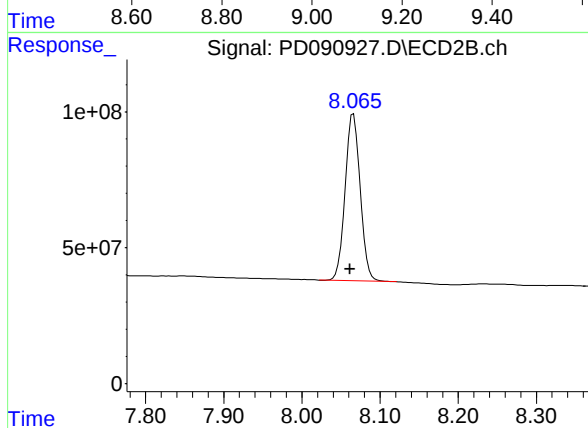
#1 Tetrachloro-m-xylene

R.T.: 2.873 min
Delta R.T.: -0.002 min
Response: 667273557
Conc: 22.41 ng/ml



#28 Decachlorobiphenyl

R.T.: 9.083 min
Delta R.T.: 0.018 min
Response: 102682607
Conc: 21.95 ng/ml



#28 Decachlorobiphenyl

R.T.: 8.066 min
Delta R.T.: 0.004 min
Response: 823925533
Conc: 27.20 ng/ml

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
 Data File : PL097705.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 13:19
 Operator : AR\AJ
 Sample : PB170340BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 ECD_L
 ClientSampleId :
 PB170340BS

A
B
C
D
E
F
G
H
I
J
K
L

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:53:49 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
 Quant Title : GC Extractables
 QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

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

System Monitoring Compounds							
1)	SA Tetrachlo...	3.524	2.821	96856817	139.3E6	23.671	23.135
28)	SA Decachlor...	9.003	7.984	69529620	105.7E6	23.370	19.769
Target Compounds							
2)	A alpha-BHC	3.971	3.325	347.0E6	494.7E6	51.068	50.846
3)	MA gamma-BHC...	4.299	3.657	325.5E6	459.5E6	50.941	50.724
4)	MA Heptachlor	4.891	4.005	327.9E6	446.8E6	51.278	49.290
5)	MB Aldrin	5.229	4.287	304.1E6	432.5E6	49.812	50.306
6)	B beta-BHC	4.487	3.954	133.7E6	188.9E6	51.312	47.817
7)	B delta-BHC	4.733	4.186	309.9E6	450.6E6	52.128	50.209
8)	B Heptachlo...	5.650	4.789	280.9E6	385.4E6	48.171	49.083
9)	A Endosulfan I	6.031	5.161	243.8E6	357.6E6	48.348	47.503
10)	B gamma-Chl...	5.904	5.042	271.5E6	415.9E6	49.530	48.054
11)	B alpha-Chl...	5.984	5.106	266.3E6	400.4E6	48.514	48.221
12)	B 4,4'-DDE	6.154	5.295	232.1E6	374.7E6	50.572	48.884
13)	MA Dieldrin	6.304	5.425	259.2E6	399.4E6	48.347	49.675
14)	MA Endrin	6.531	5.701	215.3E6	363.1E6	48.052	51.902
15)	B Endosulfa...	6.745	5.992	232.2E6	325.3E6	50.996	46.331
16)	A 4,4'-DDD	6.665	5.848	192.7E6	310.2E6	53.916	48.390
17)	MA 4,4'-DDT	6.979	6.100	197.0E6	308.0E6	48.869	44.783
18)	B Endrin al...	6.873	6.171	161.2E6	245.1E6	48.994	42.590
19)	B Endosulfa...	7.107	6.394	192.6E6	303.5E6	49.266	44.994
20)	A Methoxychlor	7.453	6.672	93021176	157.9E6	45.580	42.497
21)	B Endrin ke...	7.587	6.899	209.6E6	359.8E6	50.356	48.331
22)	Mirex	8.065	7.088	140.5E6	223.0E6	45.570	40.475

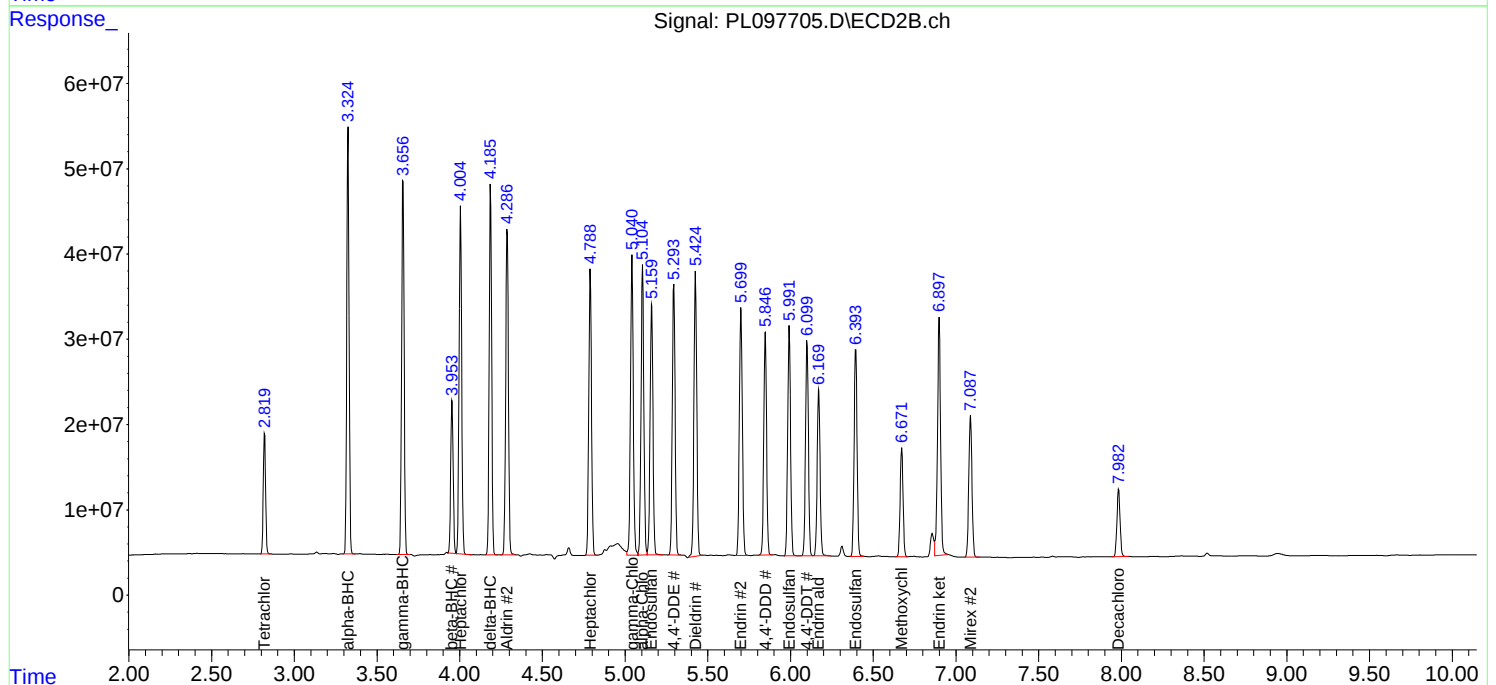
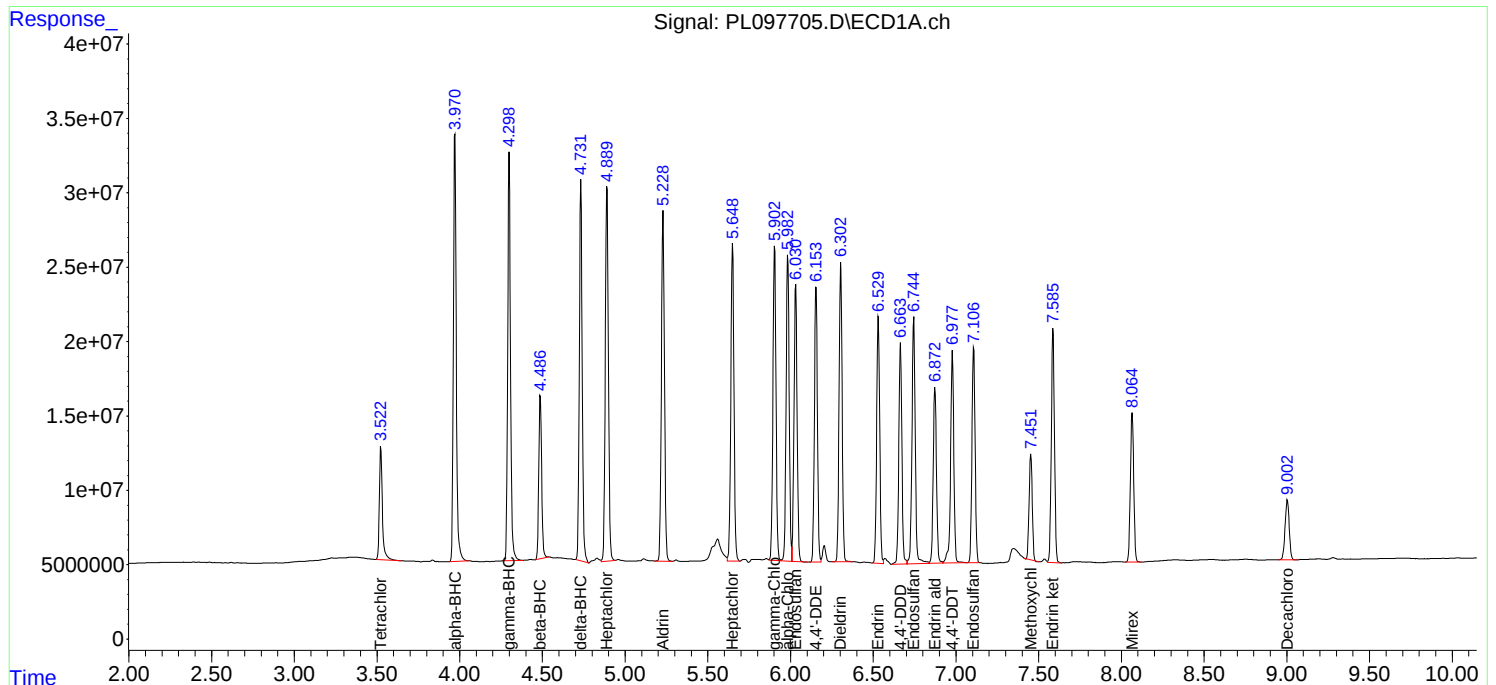
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
Data File : PL097705.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 13:19
Operator : AR\AJ
Sample : PB170340BS
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
ECD_L
ClientSampleId :
PB170340BS

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:53:49 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
Quant Title : GC Extractables
QLast Update : Tue Oct 28 17:14:31 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 x0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD110425\
 Data File : PD090952.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 04 Nov 2025 14:22
 Operator : AR\AJ
 Sample : PB170348BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PB170348BS

Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 11/05/2025

Supervised By :Yogesh Patel 11/05/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 05 00:33:16 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
 Quant Title : GC Extractables
 QLast Update : Fri Oct 17 17:43:28 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 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...	3.544	2.874	85832909	857.1E6	27.562	28.782
28)	SA Decachlor...	9.065	8.060	116.1E6	695.2E6	24.819	22.951
Target Compounds							
2)	A alpha-BHC	3.993	3.385	374.2E6	2621.2E6	52.133	51.300
3)	MA gamma-BHC...	4.324	3.721	349.4E6	2412.4E6	51.384	51.058
4)	MA Heptachlor	4.922	4.073	352.9E6	2390.9E6	63.145	53.605
5)	MB Aldrin	5.263	4.359	336.2E6	2341.8E6	51.342	50.558
6)	B beta-BHC	4.511	4.018	130.0E6	993.4E6	49.627	50.095
7)	B delta-BHC	4.759	4.253	346.9E6	2421.6E6	51.307	50.750
8)	B Heptachlo...	5.683	4.862	290.9E6	2128.4E6	50.253	49.918
9)	A Endosulfan I	6.067	5.236	270.8E6	1982.4E6	48.911	50.328
10)	B gamma-Chl...	5.939	5.115	296.6E6	2261.6E6	50.011	50.315
11)	B alpha-Chl...	6.018	5.180	299.9E6	2158.2E6	48.523m	50.031
12)	B 4,4'-DDE	6.189	5.364	255.3E6	2179.3E6	47.923	49.711
13)	MA Dieldrin	6.340	5.502	285.2E6	2183.2E6	48.904	49.412
14)	MA Endrin	6.567	5.778	239.8E6	1934.9E6	48.418	47.782
15)	B Endosulfa...	6.780	6.070	242.0E6	1767.0E6	48.402	47.051
16)	A 4,4'-DDD	6.699	5.918	201.3E6	1788.6E6	49.660	49.087
17)	MA 4,4'-DDT	7.015	6.172	197.1E6	1649.0E6	47.949	47.017
18)	B Endrin al...	6.909	6.248	164.1E6	1245.3E6	44.814	44.970
19)	B Endosulfa...	7.144	6.472	213.2E6	1657.6E6	45.964	46.243
20)	A Methoxychlor	7.488	6.743	102.3E6	788.2E6	48.118	45.225
21)	B Endrin ke...	7.625	6.981	223.4E6	1652.0E6	45.978	43.799
22)	Mirex	8.107	7.173	138.7E6	1057.5E6	43.330	41.164

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

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD110425\
Data File : PD090952.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 04 Nov 2025 14:22
Operator : AR\AJ
Sample : PB170348BS
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :

ECD_D

ClientSampleId :

PB170348BS

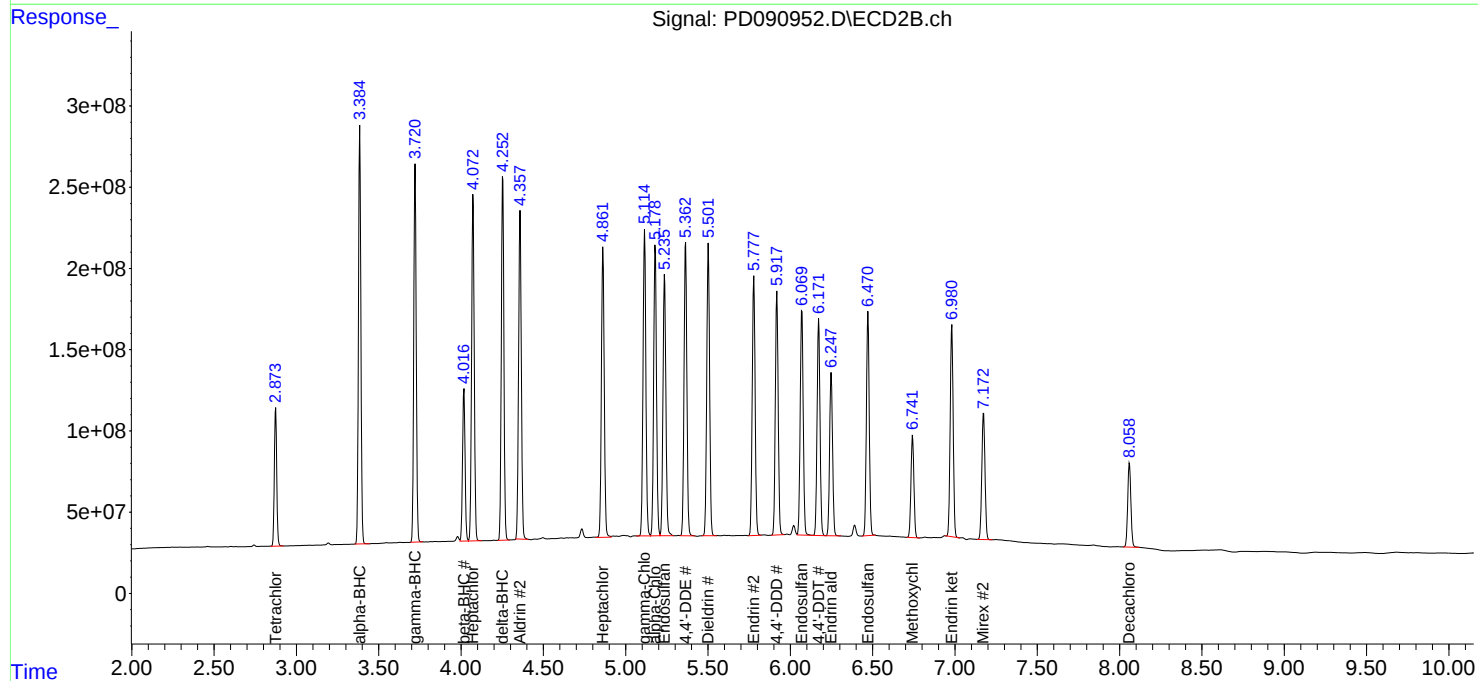
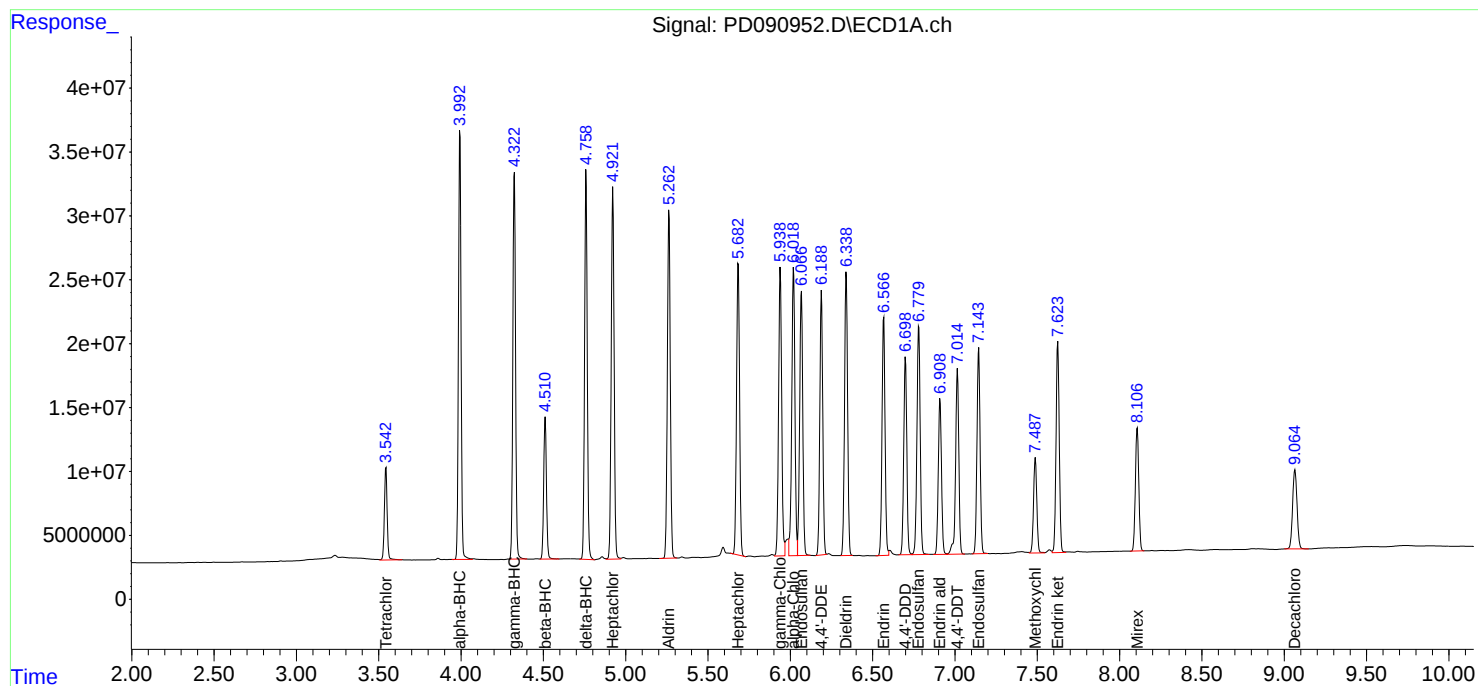
Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 11/05/2025
Supervised By :Yogesh Patel 11/05/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 05 00:33:16 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
Quant Title : GC Extractables
QLast Update : Fri Oct 17 17:43:28 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 x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD110425\
 Data File : PD090950.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 04 Nov 2025 13:52
 Operator : AR\AJ
 Sample : PB170348BSD
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 ECD_D
 ClientSampleId :
 PB170348BSD

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 05 00:32:33 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
 Quant Title : GC Extractables
 QLast Update : Fri Oct 17 17:43:28 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 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	2.873	83592138	848.1E6	26.843	28.481
28)	SA Decachlor...	9.073	8.062	116.9E6	644.4E6	24.986	21.275
Target Compounds							
2)	A alpha-BHC	4.001	3.385	373.7E6	2609.7E6	52.072	51.074
3)	MA gamma-BHC...	4.332	3.722	349.7E6	2392.7E6	51.433	50.641
4)	MA Heptachlor	4.930	4.074	354.5E6	2372.7E6	63.426	53.197
5)	MB Aldrin	5.271	4.360	334.4E6	2334.5E6	51.082	50.401
6)	B beta-BHC	4.519	4.018	127.7E6	986.4E6	48.743	49.744
7)	B delta-BHC	4.767	4.254	346.7E6	2408.5E6	51.274	50.474
8)	B Heptachlo...	5.691	4.864	281.8E6	2122.9E6	48.672	49.788
9)	A Endosulfan I	6.075	5.237	268.5E6	1933.9E6	48.490	49.095
10)	B gamma-Chl...	5.947	5.116	287.1E6	2216.7E6	48.397	49.316
11)	B alpha-Chl...	6.027	5.181	288.3E6	2109.1E6	46.642	48.893
12)	B 4,4'-DDE	6.197	5.365	257.0E6	2107.6E6	48.248	48.076
13)	MA Dieldrin	6.348	5.503	284.7E6	2124.3E6	48.829	48.079
14)	MA Endrin	6.575	5.780	235.1E6	1857.5E6	47.461	45.871
15)	B Endosulfa...	6.788	6.072	238.4E6	1699.9E6	47.683	45.263
16)	A 4,4'-DDD	6.707	5.920	198.7E6	1718.2E6	49.003	47.154
17)	MA 4,4'-DDT	7.022	6.174	195.3E6	1612.9E6	47.504	45.987
18)	B Endrin al...	6.917	6.249	162.7E6	1232.7E6	44.425	44.514
19)	B Endosulfa...	7.152	6.473	210.1E6	1570.9E6	45.291	43.826
20)	A Methoxychlor	7.495	6.744	99753193	757.8E6	46.930	43.484
21)	B Endrin ke...	7.632	6.983	222.1E6	1555.4E6	45.719	41.239
22)	Mirex	8.114	7.175	137.6E6	979.0E6	42.992	38.106

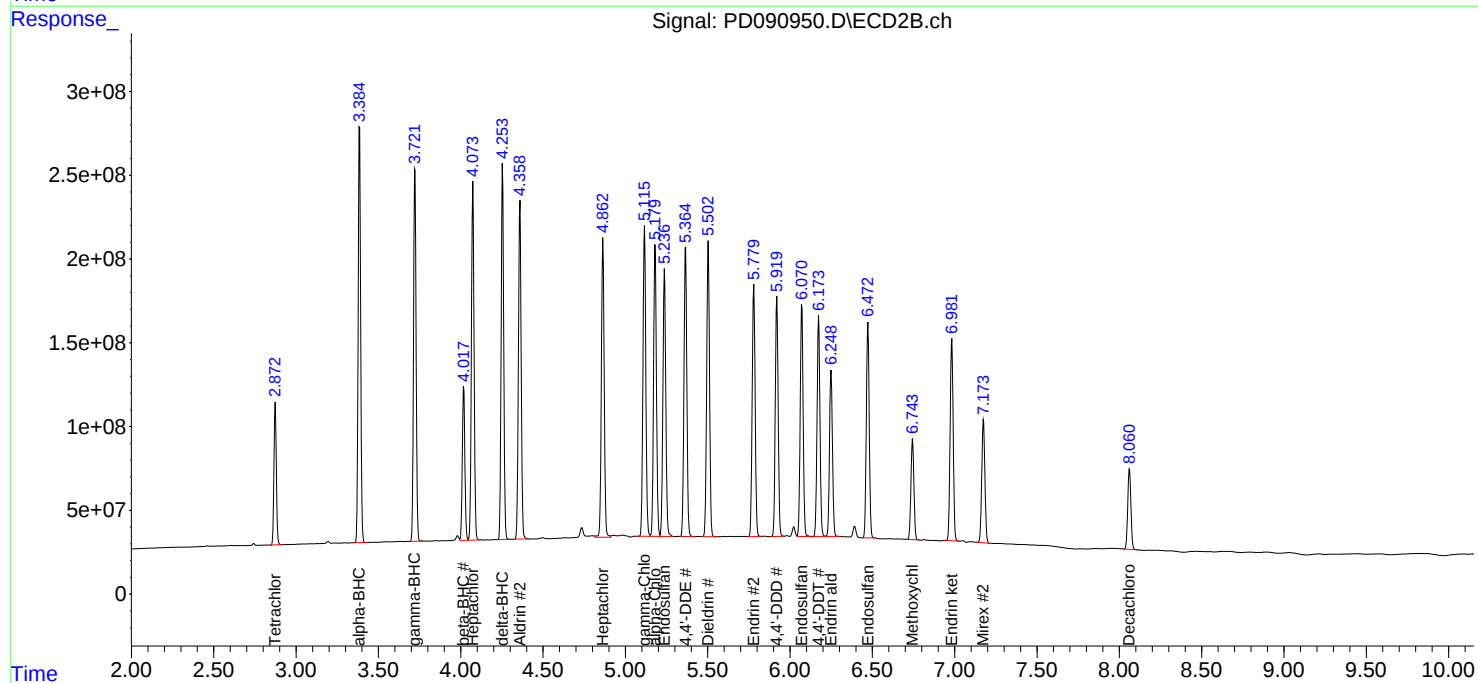
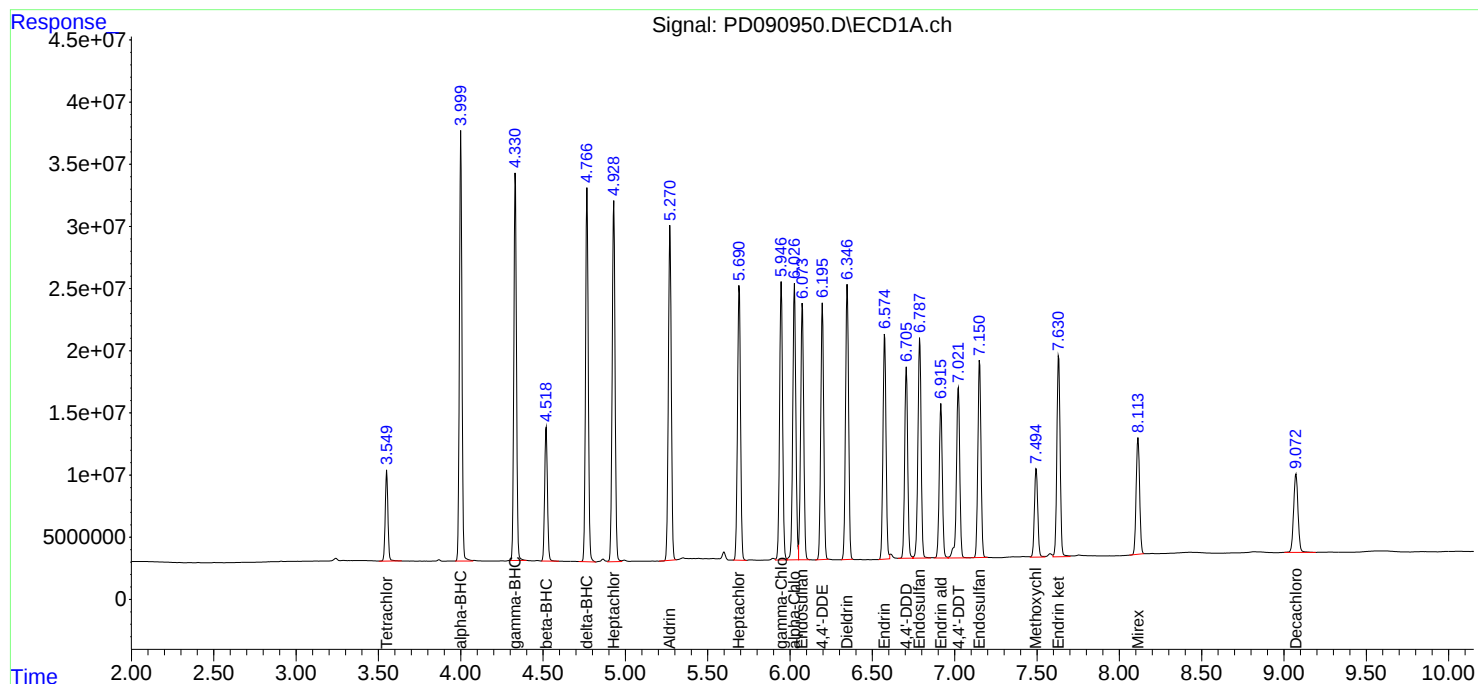
(f)=RT Delta > 1/2 Window (#)=Amounts differ by > 25% (m)=manual int.

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_D\Data\PD110425\
Data File : PD090950.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 04 Nov 2025 13:52
Operator : AR\AJ
Sample : PB170348BSD
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
ECD_D
ClientSampleId :
PB170348BSD

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 05 00:32:33 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_D\Method\PD101725.M
Quant Title : GC Extractables
QLast Update : Fri Oct 17 17:43:28 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 x0.5 Signal #2 Info : 30M x 0.32mm x 0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
 Data File : PL097707.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 13:47
 Operator : AR\AJ
 Sample : Q3495-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :

ECD_L

ClientSampleId :

SB-1025-2MS

Manual Integrations**APPROVED**

Reviewed By :Rahul Chavli 11/03/2025

Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:54:17 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
 Quant Title : GC Extractables
 QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

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

System Monitoring Compounds							
1)	SA Tetrachlo...	3.523	2.822	68747010	100.5E6	16.801m	16.695
28)	SA Decachlor...	9.001	7.984	43898537	68422749	14.755	12.802
Target Compounds							
2)	A alpha-BHC	3.972	3.327	338.6E6	477.3E6	49.836	49.061
3)	MA gamma-BHC...	4.300	3.659	318.2E6	446.5E6	49.806	49.285
4)	MA Heptachlor	4.891	4.007	331.0E6	436.9E6	51.753	48.189
5)	MB Aldrin	5.230	4.289	299.5E6	423.7E6	49.053	49.286
6)	B beta-BHC	4.488	3.956	132.4E6	184.9E6	50.820	46.806
7)	B delta-BHC	4.733	4.188	308.2E6	439.4E6	51.844	48.964
8)	B Heptachlo...	5.651	4.791	277.6E6	380.5E6	47.607	48.465
9)	A Endosulfan I	6.032	5.162	243.9E6	359.9E6	48.371	47.806
10)	B gamma-Chl...	5.904	5.043	270.6E6	412.2E6	49.358	47.624
11)	B alpha-Chl...	5.984	5.107	265.1E6	398.7E6	48.285	48.015
12)	B 4,4'-DDE	6.155	5.296	233.9E6	372.3E6	50.959	48.568
13)	MA Dieldrin	6.304	5.427	260.5E6	390.9E6	48.578	48.628
14)	MA Endrin	6.531	5.701	214.3E6	348.4E6	47.839	49.801
15)	B Endosulfa...	6.744	5.994	225.0E6	332.2E6	49.409	47.301
16)	A 4,4'-DDD	6.664	5.849	187.9E6	293.1E6	52.561	45.715
17)	MA 4,4'-DDT	6.979	6.101	201.1E6	319.8E6	49.867	46.489
18)	B Endrin al...	6.873	6.172	160.4E6	249.4E6	48.759	43.335
19)	B Endosulfa...	7.107	6.395	193.2E6	308.0E6	49.413	45.648
20)	A Methoxychlor	7.452	6.673	99197682	164.0E6	48.607	44.137
21)	B Endrin ke...	7.587	6.898	203.3E6	334.0E6	48.848	44.867m
22)	Mirex	8.065	7.090	139.3E6	229.2E6	45.165	41.592

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

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
Data File : PL097707.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 13:47
Operator : AR\AJ
Sample : Q3495-02MS
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :

ECD_L

ClientSampleId :

SB-1025-2MS

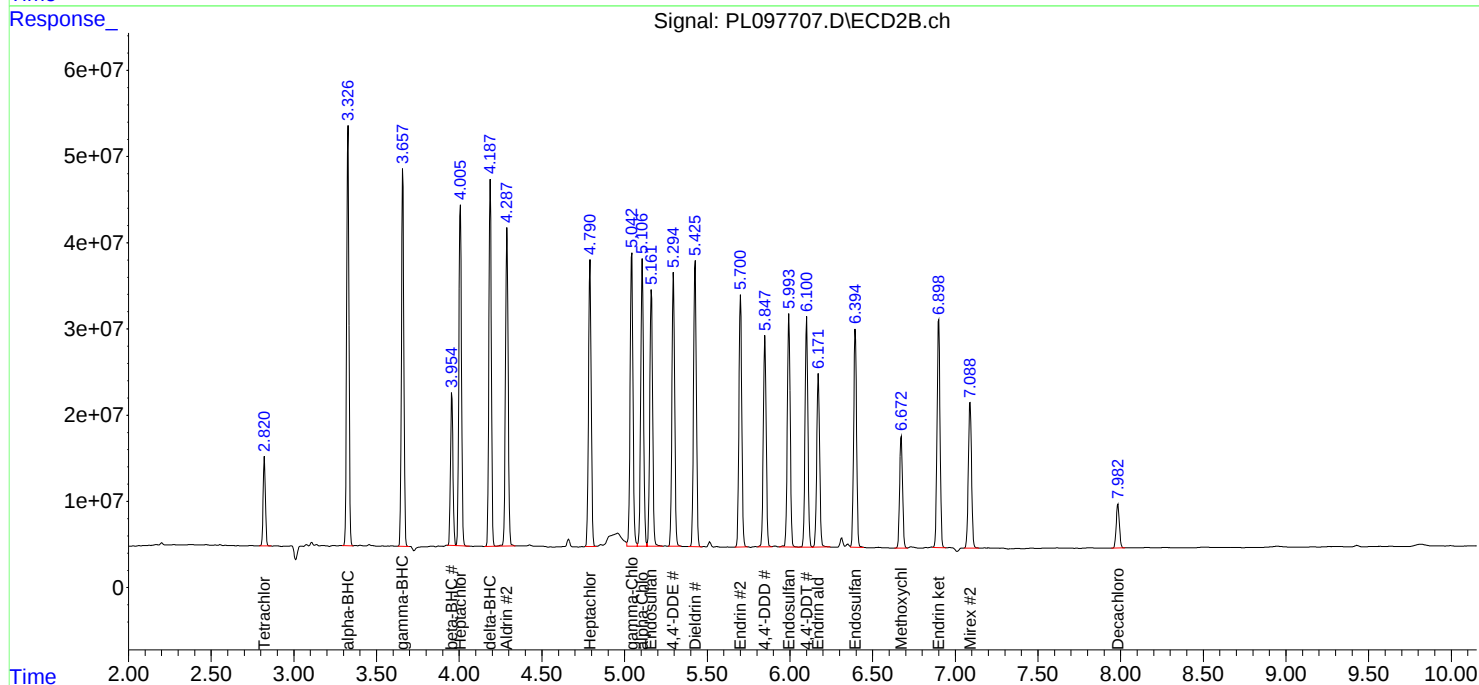
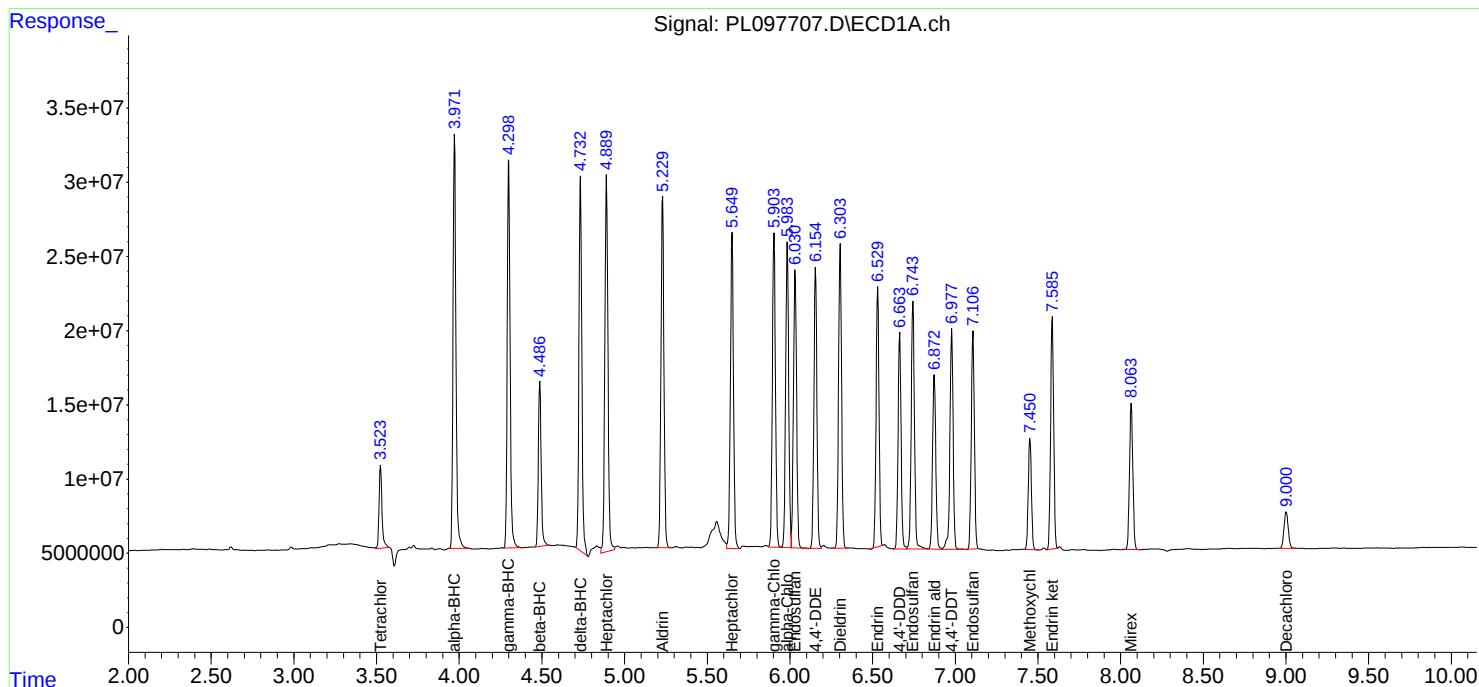
Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 11/03/2025
Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:54:17 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
Quant Title : GC Extractables
QLast Update : Tue Oct 28 17:14:31 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 x0.25µm



Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
 Data File : PL097708.D
 Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
 Acq On : 31 Oct 2025 14:00
 Operator : AR\AJ
 Sample : Q3495-02MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 ECD_L
 ClientSampleId :
 SB-1025-2MSD

Manual Integrations APPROVED

Reviewed By :Rahul Chavli 11/03/2025
 Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
 Integration File signal 2: autoint2.e
 Quant Time: Nov 01 01:54:33 2025
 Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
 Quant Title : GC Extractables
 QLast Update : Tue Oct 28 17:14:31 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 x0.25µm

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

System Monitoring Compounds							
1)	SA Tetrachlo...	3.523	2.822	69584258	103.4E6	17.006m	17.171
28)	SA Decachlor...	9.001	7.984	44314272	71690883	14.895	13.413
Target Compounds							
2)	A alpha-BHC	3.971	3.327	347.7E6	489.8E6	51.178	50.342
3)	MA gamma-BHC...	4.299	3.659	326.2E6	457.8E6	51.053	50.537
4)	MA Heptachlor	4.891	4.006	339.1E6	444.5E6	53.028	49.028
5)	MB Aldrin	5.230	4.289	306.0E6	433.5E6	50.115	50.418
6)	B beta-BHC	4.488	3.955	135.4E6	188.5E6	51.948	47.716
7)	B delta-BHC	4.733	4.188	316.3E6	450.5E6	53.209	50.199
8)	B Heptachlo...	5.650	4.791	284.0E6	388.4E6	48.693	49.468
9)	A Endosulfan I	6.031	5.162	250.8E6	367.9E6	49.732	48.873
10)	B gamma-Chl...	5.904	5.043	275.9E6	419.6E6	50.323	48.487
11)	B alpha-Chl...	5.984	5.107	273.0E6	406.6E6	49.737	48.969
12)	B 4,4'-DDE	6.155	5.295	241.0E6	381.0E6	52.519	49.708
13)	MA Dieldrin	6.304	5.426	265.8E6	399.5E6	49.580	49.693
14)	MA Endrin	6.531	5.701	218.7E6	355.2E6	48.813	50.766
15)	B Endosulfa...	6.745	5.993	229.7E6	340.0E6	50.444	48.417
16)	A 4,4'-DDD	6.665	5.848	191.7E6	302.5E6	53.629	47.180
17)	MA 4,4'-DDT	6.978	6.101	203.3E6	325.7E6	50.413	47.353
18)	B Endrin al...	6.873	6.172	164.7E6	255.1E6	50.056	44.326
19)	B Endosulfa...	7.107	6.395	196.5E6	316.7E6	50.264	46.951
20)	A Methoxychlor	7.452	6.673	98567291	167.8E6	48.298	45.171
21)	B Endrin ke...	7.587	6.898	209.1E6	344.3E6	50.226	46.254m
22)	Mirex	8.065	7.088	141.1E6	236.5E6	45.743	42.920

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

Data Path : Z:\pestpcbsrv\HPCHEM1\ECD_L\Data\PL103125\
Data File : PL097708.D
Signal(s) : Signal #1: ECD1A.ch Signal #2: ECD2B.ch
Acq On : 31 Oct 2025 14:00
Operator : AR\AJ
Sample : Q3495-02MSD
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :

ECD_L

ClientSampleId :

SB-1025-2MSD

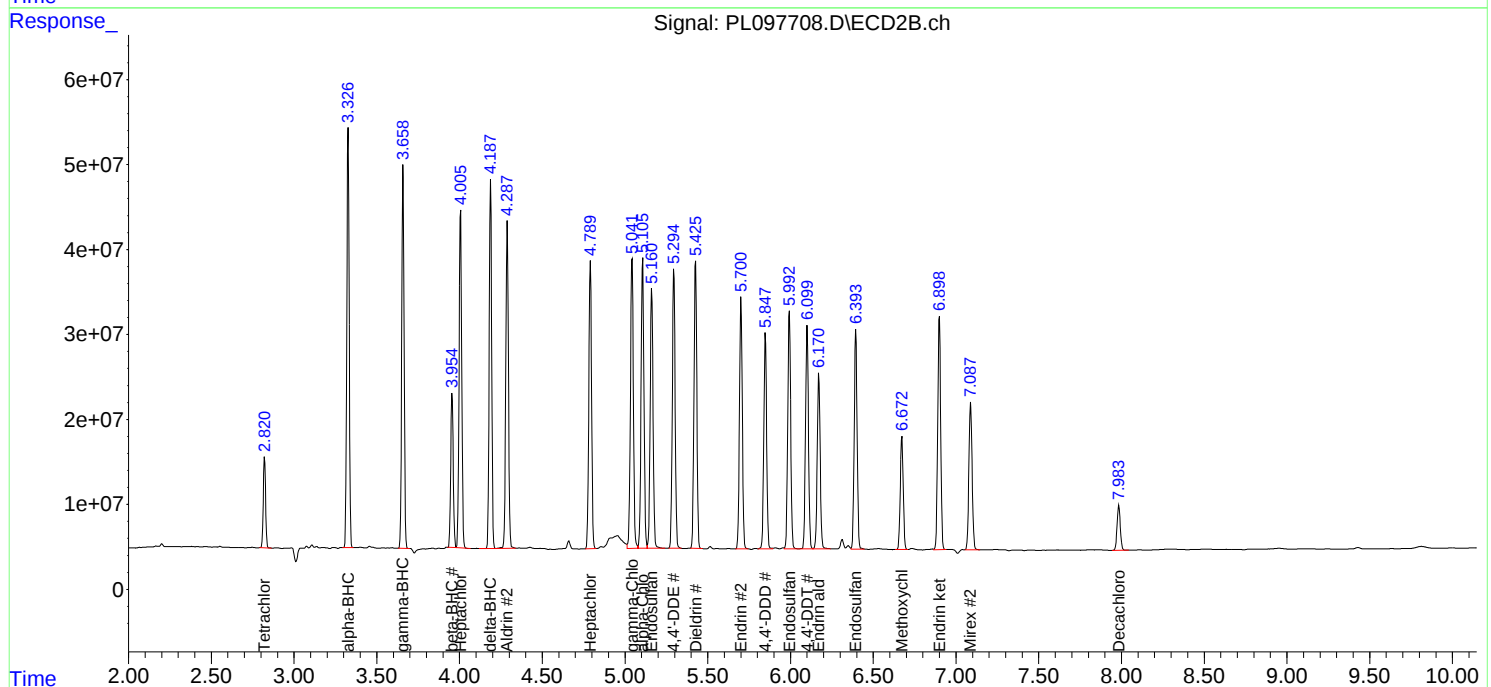
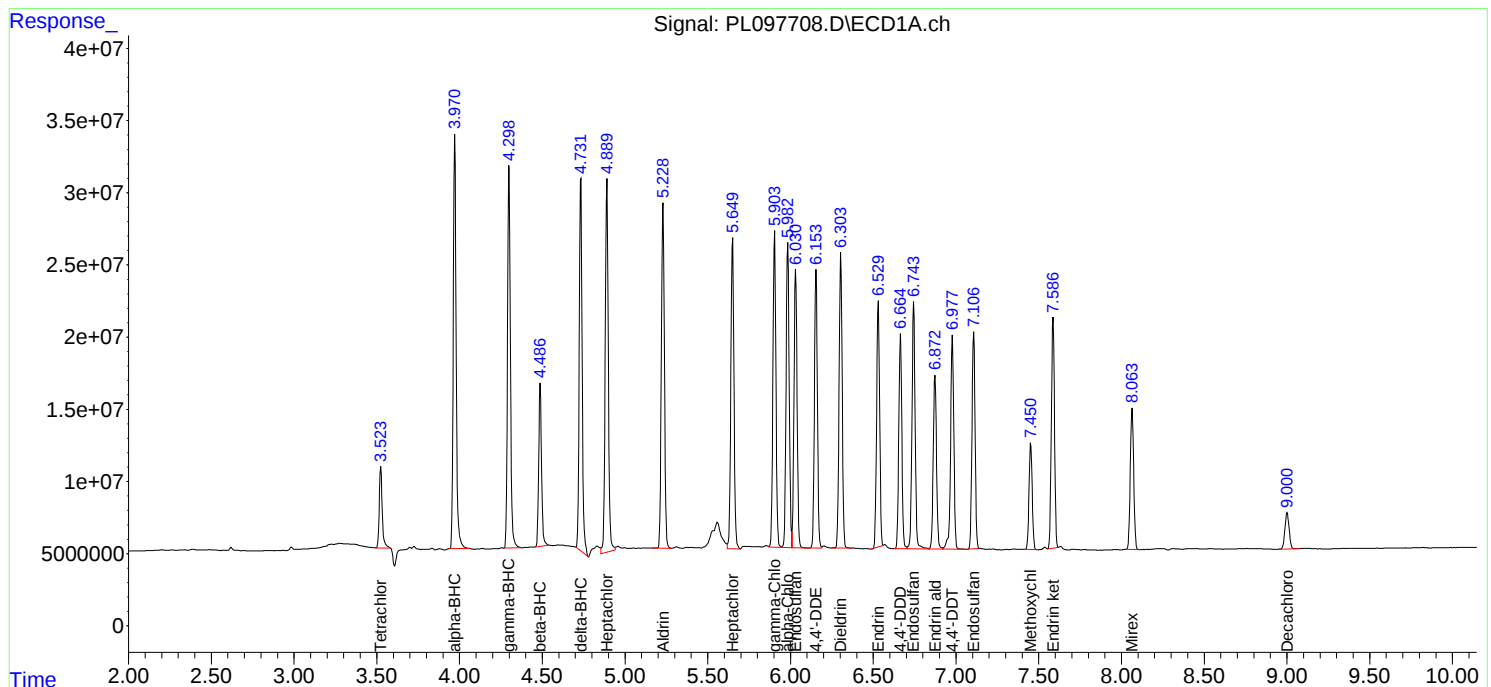
Manual Integrations

APPROVED

Reviewed By :Rahul Chavli 11/03/2025
Supervised By :Yogesh Patel 11/03/2025

Integration File signal 1: autoint1.e
Integration File signal 2: autoint2.e
Quant Time: Nov 01 01:54:33 2025
Quant Method : Z:\pestpcbsrv\HPCHEM1\ECD_L\methods\PL102825.M
Quant Title : GC Extractables
QLast Update : Tue Oct 28 17:14:31 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 x0.25µm



Manual Integration Report

Sequence:	pd101725	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PD090648.D	Endrin aldehyde	Abdul	10/19/2025 5:20:46 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PEM	PD090648.D	Endrin ketone #2	Abdul	10/19/2025 5:20:46 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
RESCHK	PD090649.D	Tetrachloro-m-xylene	Abdul	10/19/2025 5:20:37 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PSTDICC005	PD090654.D	alpha-Chlordane	Abdul	10/19/2025 5:19:40 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PSTDICC005	PD090654.D	Endrin ketone	Abdul	10/19/2025 5:19:40 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PSTDICC005	PD090654.D	Tetrachloro-m-xylene	Abdul	10/19/2025 5:19:40 PM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PTOXICV500	PD090667.D	Toxaphene-5 #2	Abdul	10/21/2025 8:49:16 AM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PEM	PD090669.D	4,4"-DDD	Abdul	10/21/2025 8:49:12 AM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PEM	PD090669.D	4,4"-DDD #2	Abdul	10/21/2025 8:49:12 AM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software
PEM	PD090669.D	Endrin aldehyde	Abdul	10/21/2025 8:49:12 AM	mohammad	10/24/2025 1:04:58	Peak Integrated by Software

Manual Integration Report

Sequence:	pd103125	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PD090925.D	Endrin aldehyde	rahul	11/3/2025 8:18:48 AM	yogesh	11/3/2025 8:49:33	Peak Integrated by Software
PSTDCCC050	PD090926.D	Endrin ketone #2	rahul	11/3/2025 8:18:51 AM	yogesh	11/3/2025 8:49:35	Peak Integrated by Software
Q3495-03	PD090931.D	Tetrachloro-m-xylene	rahul	11/3/2025 8:18:59 AM	yogesh	11/3/2025 8:49:42	Peak Integrated by Software
PSTDCCC050	PD090933.D	Endrin ketone #2	rahul	11/3/2025 8:19:03 AM	yogesh	11/3/2025 8:49:43	Peak Integrated by Software

Manual Integration Report

Sequence:	Pd110425	Instrument	ECD_d
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PSTDCCC050	PD090937.D	Endrin ketone #2	rahul	11/5/2025 8:29:24 AM	yogesh	11/5/2025 8:39:44	Peak Integrated by Software
PB170348BS	PD090952.D	alpha-Chlordane	rahul	11/5/2025 8:29:59 AM	yogesh	11/5/2025 8:40:08	Peak Integrated by Software

Manual Integration Report

Sequence:	PL102825	Instrument	ECD_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PL097637.D	4,4"-DDD	Abdul	10/29/2025 8:11:45 AM	Sohil	10/29/2025 4:11:33	Peak Integrated by Software
PEM	PL097637.D	4,4"-DDD #2	Abdul	10/29/2025 8:11:45 AM	Sohil	10/29/2025 4:11:33	Peak Integrated by Software
PEM	PL097637.D	Endrin aldehyde	Abdul	10/29/2025 8:11:45 AM	Sohil	10/29/2025 4:11:33	Peak Integrated by Software
PEM	PL097637.D	Endrin ketone	Abdul	10/29/2025 8:11:45 AM	Sohil	10/29/2025 4:11:33	Peak Integrated by Software
PEM	PL097637.D	Endrin ketone #2	Abdul	10/29/2025 8:11:45 AM	Sohil	10/29/2025 4:11:33	Peak Integrated by Software
RESCHK	PL097638.D	gamma-Chlordane #2	Abdul	10/29/2025 8:11:40 AM	Sohil	10/29/2025 4:11:35	Peak Integrated by Software
PSTDICC025	PL097642.D	Endosulfan II	Abdul	10/29/2025 8:11:33 AM	Sohil	10/29/2025 4:11:37	Peak Integrated by Software
PSTDICC005	PL097643.D	4,4"-DDD	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	4,4"-DDD #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	alpha-Chlordane #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	Endosulfan I #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	Endrin	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	Endrin #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software

Manual Integration Report

Sequence:	PL102825	Instrument	ECD_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PSTDICC005	PL097643.D	Endrin aldehyde #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	gamma-Chlordane #2	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PSTDICC005	PL097643.D	Heptachlor	Abdul	10/29/2025 8:11:28 AM	Sohil	10/29/2025 4:11:39	Peak Integrated by Software
PCHLORICV50 0	PL097655.D	Chlordane-1	Abdul	10/29/2025 8:11:21 AM	Sohil	10/29/2025 4:11:43	Peak Integrated by Software
PCHLORICV50 0	PL097655.D	Decachlorobiphenyl	Abdul	10/29/2025 8:11:21 AM	Sohil	10/29/2025 4:11:43	Peak Integrated by Software
PEM	PL097658.D	4,4"-DDD	Abdul	10/29/2025 8:11:07 AM	Sohil	10/29/2025 4:11:45	Peak Integrated by Software
PEM	PL097658.D	4,4"-DDD #2	Abdul	10/29/2025 8:11:07 AM	Sohil	10/29/2025 4:11:45	Peak Integrated by Software
PEM	PL097658.D	Endrin aldehyde	Abdul	10/29/2025 8:11:07 AM	Sohil	10/29/2025 4:11:45	Peak Integrated by Software
PEM	PL097658.D	Endrin aldehyde #2	Abdul	10/29/2025 8:11:07 AM	Sohil	10/29/2025 4:11:45	Peak Integrated by Software
PEM	PL097658.D	Endrin ketone	Abdul	10/29/2025 8:11:07 AM	Sohil	10/29/2025 4:11:45	Peak Integrated by Software

Manual Integration Report

Sequence:	pl103125	Instrument	ECD_I
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Sample ID	File ID	Parameter	Review By	Review On	Supervised By	Supervised On	Reason
PEM	PL097700.D	4,4"-DDD	rahul	11/3/2025 8:09:14 AM	yogesh	11/3/2025 8:42:42	Peak Integrated by Software
PEM	PL097700.D	Endrin aldehyde	rahul	11/3/2025 8:09:14 AM	yogesh	11/3/2025 8:42:42	Peak Integrated by Software
PSTDCCC050	PL097701.D	4,4"-DDD	rahul	11/3/2025 8:09:17 AM	yogesh	11/3/2025 8:42:45	Peak Integrated by Software
PSTDCCC050	PL097701.D	Dieldrin #2	rahul	11/3/2025 8:09:17 AM	yogesh	11/3/2025 8:42:45	Peak Integrated by Software
PSTDCCC050	PL097701.D	Endrin	rahul	11/3/2025 8:09:17 AM	yogesh	11/3/2025 8:42:45	Peak Integrated by Software
PSTDCCC050	PL097701.D	gamma-Chlordane	rahul	11/3/2025 8:09:17 AM	yogesh	11/3/2025 8:42:45	Peak Integrated by Software
PSTDCCC050	PL097701.D	Heptachlor	rahul	11/3/2025 8:09:17 AM	yogesh	11/3/2025 8:42:45	Peak Integrated by Software
Q3495-02	PL097706.D	Tetrachloro-m-xylene	rahul	11/3/2025 8:09:23 AM	yogesh	11/3/2025 8:42:53	Peak Integrated by Software
Q3495-02MS	PL097707.D	Endrin ketone #2	rahul	11/3/2025 8:09:59 AM	yogesh	11/3/2025 8:42:55	Peak Integrated by Software
Q3495-02MS	PL097707.D	Tetrachloro-m-xylene	rahul	11/3/2025 8:09:59 AM	yogesh	11/3/2025 8:42:55	Peak Integrated by Software
Q3495-02MSD	PL097708.D	Endrin ketone #2	rahul	11/3/2025 8:09:28 AM	yogesh	11/3/2025 8:42:57	Peak Integrated by Software
Q3495-02MSD	PL097708.D	Tetrachloro-m-xylene	rahul	11/3/2025 8:09:28 AM	yogesh	11/3/2025 8:42:57	Peak Integrated by Software
PSTDCCC050	PL097715.D	gamma-Chlordane	rahul	11/3/2025 8:09:50 AM	yogesh	11/3/2025 8:43:25	Peak Integrated by Software



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

Manual Integration Report

Sequence:

pl103125

Instrument

ECD_I

Sample ID

File ID

Parameter

Review By

Review On

Supervised
By

Supervised On

Reason

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD101725

Review By	Abdul	Review On	10/19/2025 5:21:31 PM
Supervise By	mohammad	Supervise On	10/24/2025 1:04:58 AM
SubDirectory	PD101725	HP Acquire Method	HP Processing Method pd101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	Sampleld	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD090646.D	17 Oct 2025 10:40	AR\AJ	Ok
2	I.BLK	PD090647.D	17 Oct 2025 10:53	AR\AJ	Ok
3	PEM	PD090648.D	17 Oct 2025 11:07	AR\AJ	Ok,M
4	RESCHK	PD090649.D	17 Oct 2025 11:20	AR\AJ	Ok,M
5	PSTDICC100	PD090650.D	17 Oct 2025 11:34	AR\AJ	Ok
6	PSTDICC075	PD090651.D	17 Oct 2025 11:48	AR\AJ	Ok
7	PSTDICC050	PD090652.D	17 Oct 2025 12:01	AR\AJ	Ok
8	PSTDICC025	PD090653.D	17 Oct 2025 12:15	AR\AJ	Ok
9	PSTDICC005	PD090654.D	17 Oct 2025 12:28	AR\AJ	Ok,M
10	PCHLORICC1000	PD090655.D	17 Oct 2025 12:42	AR\AJ	Ok
11	PCHLORICC750	PD090656.D	17 Oct 2025 12:56	AR\AJ	Ok
12	PCHLORICC500	PD090657.D	17 Oct 2025 13:09	AR\AJ	Ok
13	PCHLORICC250	PD090658.D	17 Oct 2025 13:23	AR\AJ	Ok
14	PCHLORICC050	PD090659.D	17 Oct 2025 13:36	AR\AJ	Ok,M
15	PTOXICC1000	PD090660.D	17 Oct 2025 13:50	AR\AJ	Ok
16	PTOXICC750	PD090661.D	17 Oct 2025 14:03	AR\AJ	Ok
17	PTOXICC500	PD090662.D	17 Oct 2025 14:17	AR\AJ	Ok
18	PTOXICC250	PD090663.D	17 Oct 2025 14:30	AR\AJ	Ok
19	PTOXICC100	PD090664.D	17 Oct 2025 16:18	AR\AJ	Ok
20	PSTDICV050	PD090665.D	17 Oct 2025 16:46	AR\AJ	Ok
21	PCHLORICV500	PD090666.D	17 Oct 2025 17:50	AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD101725

Review By	Abdul	Review On	10/19/2025 5:21:31 PM
Supervise By	mohammad	Supervise On	10/24/2025 1:04:58 AM
SubDirectory	PD101725	HP Acquire Method	HP Processing Method pd101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

22	PTOXICV500	PD090667.D	17 Oct 2025 18:42	AR\AJ	Ok,M
23	I.BLK	PD090668.D	17 Oct 2025 19:50	AR\AJ	Ok
24	PEM	PD090669.D	17 Oct 2025 20:04	AR\AJ	Ok,M
25	PSTDCCC050	PD090670.D	17 Oct 2025 20:18	AR\AJ	Ok
26	PB170142BL	PD090671.D	17 Oct 2025 20:31	AR\AJ	Ok
27	PB170142BS	PD090672.D	17 Oct 2025 20:45	AR\AJ	Ok
28	Q3366-01	PD090673.D	17 Oct 2025 20:59	AR\AJ	Ok,M
29	Q3366-03	PD090674.D	17 Oct 2025 21:12	AR\AJ	Ok,M
30	Q3367-01	PD090675.D	17 Oct 2025 21:26	AR\AJ	Ok,M
31	Q3369-01	PD090676.D	17 Oct 2025 21:40	AR\AJ	Ok,M
32	Q3369-05	PD090677.D	17 Oct 2025 21:53	AR\AJ	Ok,M
33	Q3363-01	PD090678.D	17 Oct 2025 22:48	AR\AJ	Ok,M
34	Q3363-02	PD090679.D	17 Oct 2025 23:02	AR\AJ	Ok,M
35	Q3363-03	PD090680.D	17 Oct 2025 23:15	AR\AJ	Ok,M
36	Q3363-04	PD090681.D	17 Oct 2025 23:29	AR\AJ	Ok,M
37	Q3363-04MS	PD090682.D	17 Oct 2025 23:43	AR\AJ	Ok,M
38	Q3363-04MSD	PD090683.D	17 Oct 2025 23:56	AR\AJ	Ok,M
39	Q3363-05	PD090684.D	18 Oct 2025 00:10	AR\AJ	Ok,M
40	Q3343-04MS	PD090685.D	18 Oct 2025 00:24	AR\AJ	Ok,M
41	Q3343-04MSD	PD090686.D	18 Oct 2025 00:37	AR\AJ	Ok
42	I.BLK	PD090687.D	18 Oct 2025 00:51	AR\AJ	Ok
43	PSTDCCC050	PD090688.D	18 Oct 2025 02:13	AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD103125

Review By	rahul	Review On	11/3/2025 8:19:14 AM
Supervise By	yogesh	Supervise On	11/3/2025 8:49:55 AM
SubDirectory	PD103125	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD090923.D	31 Oct 2025 09:53	AR\AJ	Ok
2	I.BLK	PD090924.D	31 Oct 2025 10:06	AR\AJ	Ok
3	PEM	PD090925.D	31 Oct 2025 10:20	AR\AJ	Ok,M
4	PSTDCCC050	PD090926.D	31 Oct 2025 11:09	AR\AJ	Ok,M
5	PB170348BL	PD090927.D	31 Oct 2025 14:37	AR\AJ	Ok
6	PB170348BS	PD090928.D	31 Oct 2025 14:51	AR\AJ	Not Ok
7	PB170348BSD	PD090929.D	31 Oct 2025 15:07	AR\AJ	Not Ok
8	Q3477-03	PD090930.D	31 Oct 2025 15:21	AR\AJ	Ok,M
9	Q3495-03	PD090931.D	31 Oct 2025 15:34	AR\AJ	Ok,M
10	I.BLK	PD090932.D	31 Oct 2025 15:48	AR\AJ	Ok
11	PSTDCCC050	PD090933.D	31 Oct 2025 16:01	AR\AJ	Ok,M

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD110425

Review By	rahul	Review On	11/5/2025 8:30:12 AM
Supervise By	yogesh	Supervise On	11/5/2025 8:40:26 AM
SubDirectory	PD110425	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PD090934.D	04 Nov 2025 09:04	AR\AJ	Ok
2	I.BLK	PD090935.D	04 Nov 2025 09:17	AR\AJ	Ok
3	PEM	PD090936.D	04 Nov 2025 09:31	AR\AJ	Ok
4	PSTDCCC050	PD090937.D	04 Nov 2025 09:45	AR\AJ	Ok,M
5	Q3515-01	PD090938.D	04 Nov 2025 10:17	AR\AJ	Ok,M
6	Q3516-01	PD090939.D	04 Nov 2025 10:31	AR\AJ	Ok,M
7	Q3519-01	PD090940.D	04 Nov 2025 10:44	AR\AJ	Ok,M
8	Q3519-01MS	PD090941.D	04 Nov 2025 10:58	AR\AJ	Ok,M
9	Q3519-01MSD	PD090942.D	04 Nov 2025 11:12	AR\AJ	Ok,M
10	Q3520-01	PD090943.D	04 Nov 2025 11:25	AR\AJ	Ok,M
11	Q3368-06	PD090944.D	04 Nov 2025 11:39	AR\AJ	Not Ok
12	Q3368-06MS	PD090945.D	04 Nov 2025 11:53	AR\AJ	Not Ok
13	Q3368-06MSD	PD090946.D	04 Nov 2025 12:06	AR\AJ	Not Ok
14	Q3368-07	PD090947.D	04 Nov 2025 12:22	AR\AJ	Not Ok
15	I.BLK	PD090948.D	04 Nov 2025 12:35	AR\AJ	Ok
16	PSTDCCC050	PD090949.D	04 Nov 2025 13:29	AR\AJ	Ok
17	PB170348BSD	PD090950.D	04 Nov 2025 13:52	AR\AJ	Ok
18	PB170371BL	PD090951.D	04 Nov 2025 14:09	AR\AJ	Ok
19	PB170348BS	PD090952.D	04 Nov 2025 14:22	AR\AJ	Ok,M
20	PB170387BS	PD090953.D	04 Nov 2025 15:41	AR\AJ	Ok,M
21	PB170387BL	PD090954.D	04 Nov 2025 16:03	AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD110425

Review By	rahul	Review On	11/5/2025 8:30:12 AM
Supervise By	yogesh	Supervise On	11/5/2025 8:40:26 AM
SubDirectory	PD110425	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

22	PB170387TB	PD090955.D	04 Nov 2025 16:17	AR\AJ	Ok
23	I.BLK	PD090956.D	04 Nov 2025 16:31	AR\AJ	Ok
24	PSTDCCC050	PD090957.D	04 Nov 2025 16:44	AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL102825

Review By	Abdul	Review On	10/29/2025 8:12:28 AM
Supervise By	Sohil	Supervise On	10/29/2025 4:12:00 PM
SubDirectory	PL102825	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PL097635.D	28 Oct 2025 14:03	AR\AJ	Ok
2	I.BLK	PL097636.D	28 Oct 2025 14:17	AR\AJ	Ok
3	PEM	PL097637.D	28 Oct 2025 14:30	AR\AJ	Ok,M
4	RESCHK	PL097638.D	28 Oct 2025 14:44	AR\AJ	Ok,M
5	PSTDICC100	PL097639.D	28 Oct 2025 14:57	AR\AJ	Ok
6	PSTDICC075	PL097640.D	28 Oct 2025 15:11	AR\AJ	Ok
7	PSTDICC050	PL097641.D	28 Oct 2025 15:25	AR\AJ	Ok
8	PSTDICC025	PL097642.D	28 Oct 2025 15:38	AR\AJ	Ok,M
9	PSTDICC005	PL097643.D	28 Oct 2025 15:52	AR\AJ	Ok,M
10	PCHLORICC1000	PL097644.D	28 Oct 2025 16:05	AR\AJ	Ok
11	PCHLORICC750	PL097645.D	28 Oct 2025 16:19	AR\AJ	Ok
12	PCHLORICC500	PL097646.D	28 Oct 2025 16:32	AR\AJ	Ok
13	PCHLORICC250	PL097647.D	28 Oct 2025 16:46	AR\AJ	Ok
14	PCHLORICC050	PL097648.D	28 Oct 2025 17:00	AR\AJ	Ok,M
15	PTOXICC1000	PL097649.D	28 Oct 2025 17:13	AR\AJ	Ok
16	PTOXICC750	PL097650.D	28 Oct 2025 17:27	AR\AJ	Ok
17	PTOXICC500	PL097651.D	28 Oct 2025 17:40	AR\AJ	Ok
18	PTOXICC250	PL097652.D	28 Oct 2025 17:54	AR\AJ	Ok
19	PTOXICC100	PL097653.D	28 Oct 2025 18:08	AR\AJ	Ok,M
20	PSTDICV050	PL097654.D	28 Oct 2025 18:21	AR\AJ	Ok
21	PCHLORICV500	PL097655.D	28 Oct 2025 18:35	AR\AJ	Ok,M

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL102825

Review By	Abdul	Review On	10/29/2025 8:12:28 AM
Supervise By	Sohil	Supervise On	10/29/2025 4:12:00 PM
SubDirectory	PL102825	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

22	PTOXICV500	PL097656.D	28 Oct 2025 18:48	AR\AJ	Ok
23	I.BLK	PL097657.D	28 Oct 2025 19:02	AR\AJ	Ok
24	PEM	PL097658.D	28 Oct 2025 19:16	AR\AJ	Ok,M
25	PSTDCCC050	PL097659.D	28 Oct 2025 19:29	AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL103125

Review By	rahul	Review On	10/31/2025 2:27:02 PM
Supervise By	yogesh	Supervise On	11/3/2025 8:43:32 AM
SubDirectory	PL103125	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	Data File Name	Date-Time	Operator	Status
1	HEXANE	PL097698.D	31 Oct 2025 09:28	AR\AJ	Ok
2	I.BLK	PL097699.D	31 Oct 2025 09:42	AR\AJ	Ok
3	PEM	PL097700.D	31 Oct 2025 09:56	AR\AJ	Ok,M
4	PSTDCCC050	PL097701.D	31 Oct 2025 11:02	AR\AJ	Ok,M
5	Q3488-02	PL097702.D	31 Oct 2025 12:03	AR\AJ	Not Ok
6	Q3490-01	PL097703.D	31 Oct 2025 12:17	AR\AJ	Ok,M
7	PB170340BL	PL097704.D	31 Oct 2025 13:06	AR\AJ	Ok
8	PB170340BS	PL097705.D	31 Oct 2025 13:19	AR\AJ	Ok
9	Q3495-02	PL097706.D	31 Oct 2025 13:33	AR\AJ	Ok,M
10	Q3495-02MS	PL097707.D	31 Oct 2025 13:47	AR\AJ	Ok,M
11	Q3495-02MSD	PL097708.D	31 Oct 2025 14:00	AR\AJ	Ok,M
12	Q3496-01	PL097709.D	31 Oct 2025 14:14	AR\AJ	Ok,M
13	Q3497-01	PL097710.D	31 Oct 2025 14:27	AR\AJ	Ok,M
14	Q3500-01	PL097711.D	31 Oct 2025 14:41	AR\AJ	Ok,M
15	Q3501-01	PL097712.D	31 Oct 2025 14:55	AR\AJ	Ok,M
16	Q3502-01	PL097713.D	31 Oct 2025 15:08	AR\AJ	Ok,M
17	I.BLK	PL097714.D	31 Oct 2025 15:49	AR\AJ	Ok
18	PSTDCCC050	PL097715.D	31 Oct 2025 16:03	AR\AJ	Ok,M

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD101725

Review By	Abdul	Review On	10/19/2025 5:21:31 PM
Supervise By	mohammad	Supervise On	10/24/2025 1:04:58 AM
SubDirectory	PD101725	HP Acquire Method	HP Processing Method pd101725 8081

STD. NAME	STD REF.#
Tune/Reschk	PP24942,PP24651
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P P24763,PP24764
CCC	PP24751,PP24756,PP24761
Internal Standard/PEM	
ICV/I.BLK	PP24754,PP24759,PP24761
Surrogate Standard	
MS/MSD Standard	
LCS Standard	

Sr#	Sampleld	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD090646.D	17 Oct 2025 10:40		AR\AJ	Ok
2	I.BLK	I.BLK	PD090647.D	17 Oct 2025 10:53		AR\AJ	Ok
3	PEM	PEM	PD090648.D	17 Oct 2025 11:07		AR\AJ	Ok,M
4	RESCHK	RESCHK	PD090649.D	17 Oct 2025 11:20		AR\AJ	Ok,M
5	PSTDICC100	PSTDICC100	PD090650.D	17 Oct 2025 11:34		AR\AJ	Ok
6	PSTDICC075	PSTDICC075	PD090651.D	17 Oct 2025 11:48		AR\AJ	Ok
7	PSTDICC050	PSTDICC050	PD090652.D	17 Oct 2025 12:01		AR\AJ	Ok
8	PSTDICC025	PSTDICC025	PD090653.D	17 Oct 2025 12:15		AR\AJ	Ok
9	PSTDICC005	PSTDICC005	PD090654.D	17 Oct 2025 12:28		AR\AJ	Ok,M
10	PCHLORICC1000	PCHLORICC1000	PD090655.D	17 Oct 2025 12:42		AR\AJ	Ok
11	PCHLORICC750	PCHLORICC750	PD090656.D	17 Oct 2025 12:56		AR\AJ	Ok
12	PCHLORICC500	PCHLORICC500	PD090657.D	17 Oct 2025 13:09		AR\AJ	Ok
13	PCHLORICC250	PCHLORICC250	PD090658.D	17 Oct 2025 13:23		AR\AJ	Ok
14	PCHLORICC050	PCHLORICC050	PD090659.D	17 Oct 2025 13:36		AR\AJ	Ok,M
15	PTOXICC1000	PTOXICC1000	PD090660.D	17 Oct 2025 13:50		AR\AJ	Ok
16	PTOXICC750	PTOXICC750	PD090661.D	17 Oct 2025 14:03		AR\AJ	Ok
17	PTOXICC500	PTOXICC500	PD090662.D	17 Oct 2025 14:17		AR\AJ	Ok
18	PTOXICC250	PTOXICC250	PD090663.D	17 Oct 2025 14:30		AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD101725

Review By	Abdul	Review On	10/19/2025 5:21:31 PM
Supervise By	mohammad	Supervise On	10/24/2025 1:04:58 AM
SubDirectory	PD101725	HP Acquire Method	HP Processing Method pd101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

19	PTOXICC100	PTOXICC100	PD090664.D	17 Oct 2025 16:18		AR\AJ	Ok
20	PSTDICV050	ICVPD101725	PD090665.D	17 Oct 2025 16:46		AR\AJ	Ok
21	PCHLORICV500	ICVPD101725CHLOR	PD090666.D	17 Oct 2025 17:50		AR\AJ	Ok
22	PTOXICV500	ICVPD101725TOX	PD090667.D	17 Oct 2025 18:42		AR\AJ	Ok,M
23	I.BLK	I.BLK	PD090668.D	17 Oct 2025 19:50		AR\AJ	Ok
24	PEM	PEM	PD090669.D	17 Oct 2025 20:04		AR\AJ	Ok,M
25	PSTDCCC050	PSTDCCC050	PD090670.D	17 Oct 2025 20:18		AR\AJ	Ok
26	PB170142BL	PB170142BL	PD090671.D	17 Oct 2025 20:31		AR\AJ	Ok
27	PB170142BS	PB170142BS	PD090672.D	17 Oct 2025 20:45		AR\AJ	Ok
28	Q3366-01	TR-06-101625	PD090673.D	17 Oct 2025 20:59		AR\AJ	Ok,M
29	Q3366-03	TR-1-101625	PD090674.D	17 Oct 2025 21:12		AR\AJ	Ok,M
30	Q3367-01	BU-03-101625	PD090675.D	17 Oct 2025 21:26		AR\AJ	Ok,M
31	Q3369-01	MH-22	PD090676.D	17 Oct 2025 21:40		AR\AJ	Ok,M
32	Q3369-05	MH-21	PD090677.D	17 Oct 2025 21:53		AR\AJ	Ok,M
33	Q3363-01	PR132-SO1-085011-20	PD090678.D	17 Oct 2025 22:48		AR\AJ	Ok,M
34	Q3363-02	PR132-SO1-039085-20	PD090679.D	17 Oct 2025 23:02		AR\AJ	Ok,M
35	Q3363-03	PR132-SO3-028010-20	PD090680.D	17 Oct 2025 23:15		AR\AJ	Ok,M
36	Q3363-04	PR132-SO3-010013-20	PD090681.D	17 Oct 2025 23:29		AR\AJ	Ok,M
37	Q3363-04MS	PR132-SO3-010013-20	PD090682.D	17 Oct 2025 23:43		AR\AJ	Ok,M

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD101725

Review By	Abdul	Review On	10/19/2025 5:21:31 PM
Supervise By	mohammad	Supervise On	10/24/2025 1:04:58 AM
SubDirectory	PD101725	HP Acquire Method	HP Processing Method pd101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

38	Q3363-04MSD	PR132-SO3-010013-20	PD090683.D	17 Oct 2025 23:56		AR\AJ	Ok,M
39	Q3363-05	COMP01-20251015	PD090684.D	18 Oct 2025 00:10		AR\AJ	Ok,M
40	Q3343-04MS	SOIL-ROLLOFF-WC-1	PD090685.D	18 Oct 2025 00:24		AR\AJ	Ok,M
41	Q3343-04MSD	SOIL-ROLLOFF-WC-1	PD090686.D	18 Oct 2025 00:37		AR\AJ	Ok
42	I.BLK	I.BLK	PD090687.D	18 Oct 2025 00:51		AR\AJ	Ok
43	PSTDCCC050	PSTDCCC050	PD090688.D	18 Oct 2025 02:13		AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD103125

Review By	rahul	Review On	11/3/2025 8:19:14 AM
Supervise By	yogesh	Supervise On	11/3/2025 8:49:55 AM
SubDirectory	PD103125	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD090923.D	31 Oct 2025 09:53		AR\AJ	Ok
2	I.BLK	I.BLK	PD090924.D	31 Oct 2025 10:06		AR\AJ	Ok
3	PEM	PEM	PD090925.D	31 Oct 2025 10:20		AR\AJ	Ok,M
4	PSTDCCC050	PSTDCCC050	PD090926.D	31 Oct 2025 11:09		AR\AJ	Ok,M
5	PB170348BL	PB170348BL	PD090927.D	31 Oct 2025 14:37		AR\AJ	Ok
6	PB170348BS	PB170348BS	PD090928.D	31 Oct 2025 14:51	recovery fail	AR\AJ	Not Ok
7	PB170348BSD	PB170348BSD	PD090929.D	31 Oct 2025 15:07	recovery fail	AR\AJ	Not Ok
8	Q3477-03	SOUTH-MAHWAH-WA	PD090930.D	31 Oct 2025 15:21		AR\AJ	Ok,M
9	Q3495-03	TW-1025-1	PD090931.D	31 Oct 2025 15:34		AR\AJ	Ok,M
10	I.BLK	I.BLK	PD090932.D	31 Oct 2025 15:48		AR\AJ	Ok
11	PSTDCCC050	PSTDCCC050	PD090933.D	31 Oct 2025 16:01		AR\AJ	Ok,M

M : Manual Integration

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD110425

Review By	rahul	Review On	11/5/2025 8:30:12 AM
Supervise By	yogesh	Supervise On	11/5/2025 8:40:26 AM
SubDirectory	PD110425	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PD090934.D	04 Nov 2025 09:04		AR\AJ	Ok
2	I.BLK	I.BLK	PD090935.D	04 Nov 2025 09:17		AR\AJ	Ok
3	PEM	PEM	PD090936.D	04 Nov 2025 09:31		AR\AJ	Ok
4	PSTDCCC050	PSTDCCC050	PD090937.D	04 Nov 2025 09:45		AR\AJ	Ok,M
5	Q3515-01	SU-04-103125	PD090938.D	04 Nov 2025 10:17		AR\AJ	Ok,M
6	Q3516-01	HD-01-103125	PD090939.D	04 Nov 2025 10:31		AR\AJ	Ok,M
7	Q3519-01	PL-01-103125	PD090940.D	04 Nov 2025 10:44		AR\AJ	Ok,M
8	Q3519-01MS	PL-01-103125MS	PD090941.D	04 Nov 2025 10:58		AR\AJ	Ok,M
9	Q3519-01MSD	PL-01-103125MSD	PD090942.D	04 Nov 2025 11:12		AR\AJ	Ok,M
10	Q3520-01	EO-02-103125	PD090943.D	04 Nov 2025 11:25		AR\AJ	Ok,M
11	Q3368-06	OUTSIDE-DUMPSTER	PD090944.D	04 Nov 2025 11:39	All surrogate fail	AR\AJ	Not Ok
12	Q3368-06MS	OUTSIDE-DUMPSTER	PD090945.D	04 Nov 2025 11:53	DCB Low in both column , Comp#8,14 recover fail	AR\AJ	Not Ok
13	Q3368-06MSD	OUTSIDE-DUMPSTER	PD090946.D	04 Nov 2025 12:06	DCB Low in both column , Comp#8,14 recover fail	AR\AJ	Not Ok
14	Q3368-07	PRESS-SLUDGE	PD090947.D	04 Nov 2025 12:22	All surrogate fail	AR\AJ	Not Ok
15	I.BLK	I.BLK	PD090948.D	04 Nov 2025 12:35		AR\AJ	Ok
16	PSTDCCC050	PSTDCCC050	PD090949.D	04 Nov 2025 13:29	Comp#4 high in 1st column , Comp#22 and DCB low in 2nd column	AR\AJ	Ok
17	PB170348BSD	PB170348BSD	PD090950.D	04 Nov 2025 13:52		AR\AJ	Ok

Instrument ID: ECD_D

Daily Analysis Runlog For Sequence/QC Batch ID # PD110425

Review By	rahul	Review On	11/5/2025 8:30:12 AM
Supervise By	yogesh	Supervise On	11/5/2025 8:40:26 AM
SubDirectory	PD110425	HP Acquire Method	HP Processing Method PD101725 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,PP24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

18	PB170371BL	PB170371BL	PD090951.D	04 Nov 2025 14:09		AR\AJ	Ok
19	PB170348BS	PB170348BS	PD090952.D	04 Nov 2025 14:22		AR\AJ	Ok,M
20	PB170387BS	PB170387BS	PD090953.D	04 Nov 2025 15:41		AR\AJ	Ok,M
21	PB170387BL	PB170387BL	PD090954.D	04 Nov 2025 16:03		AR\AJ	Ok
22	PB170387TB	PB170387TB	PD090955.D	04 Nov 2025 16:17		AR\AJ	Ok
23	I.BLK	I.BLK	PD090956.D	04 Nov 2025 16:31		AR\AJ	Ok
24	PSTDCCC050	PSTDCCC050	PD090957.D	04 Nov 2025 16:44	Comp#4 high in 1st column	AR\AJ	Ok

M : Manual Integration

A
B
C
D
E
F
G
H
I
J
K
L

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL102825

Review By	Abdul	Review On	10/29/2025 8:12:28 AM
Supervise By	Sohil	Supervise On	10/29/2025 4:12:00 PM
SubDirectory	PL102825	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PL097635.D	28 Oct 2025 14:03		AR\AJ	Ok
2	I.BLK	I.BLK	PL097636.D	28 Oct 2025 14:17		AR\AJ	Ok
3	PEM	PEM	PL097637.D	28 Oct 2025 14:30		AR\AJ	Ok,M
4	RESCHK	RESCHK	PL097638.D	28 Oct 2025 14:44		AR\AJ	Ok,M
5	PSTDICC100	PSTDICC100	PL097639.D	28 Oct 2025 14:57		AR\AJ	Ok
6	PSTDICC075	PSTDICC075	PL097640.D	28 Oct 2025 15:11		AR\AJ	Ok
7	PSTDICC050	PSTDICC050	PL097641.D	28 Oct 2025 15:25		AR\AJ	Ok
8	PSTDICC025	PSTDICC025	PL097642.D	28 Oct 2025 15:38		AR\AJ	Ok,M
9	PSTDICC005	PSTDICC005	PL097643.D	28 Oct 2025 15:52		AR\AJ	Ok,M
10	PCHLORICC1000	PCHLORICC1000	PL097644.D	28 Oct 2025 16:05		AR\AJ	Ok
11	PCHLORICC750	PCHLORICC750	PL097645.D	28 Oct 2025 16:19		AR\AJ	Ok
12	PCHLORICC500	PCHLORICC500	PL097646.D	28 Oct 2025 16:32		AR\AJ	Ok
13	PCHLORICC250	PCHLORICC250	PL097647.D	28 Oct 2025 16:46		AR\AJ	Ok
14	PCHLORICC050	PCHLORICC050	PL097648.D	28 Oct 2025 17:00		AR\AJ	Ok,M
15	PTOXICC1000	PTOXICC1000	PL097649.D	28 Oct 2025 17:13		AR\AJ	Ok
16	PTOXICC750	PTOXICC750	PL097650.D	28 Oct 2025 17:27		AR\AJ	Ok
17	PTOXICC500	PTOXICC500	PL097651.D	28 Oct 2025 17:40		AR\AJ	Ok
18	PTOXICC250	PTOXICC250	PL097652.D	28 Oct 2025 17:54		AR\AJ	Ok

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL102825

Review By	Abdul	Review On	10/29/2025 8:12:28 AM
Supervise By	Sohil	Supervise On	10/29/2025 4:12:00 PM
SubDirectory	PL102825	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

19	PTOXICC100	PTOXICC100	PL097653.D	28 Oct 2025 18:08		AR\AJ	Ok,M
20	PSTDICV050	ICVPL102825	PL097654.D	28 Oct 2025 18:21		AR\AJ	Ok
21	PCHLORICV500	ICVPL102825CHLOR	PL097655.D	28 Oct 2025 18:35		AR\AJ	Ok,M
22	PTOXICV500	ICVPL102825TOX	PL097656.D	28 Oct 2025 18:48		AR\AJ	Ok
23	I.BLK	I.BLK	PL097657.D	28 Oct 2025 19:02		AR\AJ	Ok
24	PEM	PEM	PL097658.D	28 Oct 2025 19:16		AR\AJ	Ok,M
25	PSTDCCC050	PSTDCCC050	PL097659.D	28 Oct 2025 19:29		AR\AJ	Ok

M : Manual Integration

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL103125

Review By	rahul	Review On	10/31/2025 2:27:02 PM
Supervise By	yogesh	Supervise On	11/3/2025 8:43:32 AM
SubDirectory	PL103125	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

Sr#	SampleId	ClientID	Data File Name	Date-Time	Comment	Operator	Status
1	HEXANE	HEXANE	PL097698.D	31 Oct 2025 09:28		AR\AJ	Ok
2	I.BLK	I.BLK	PL097699.D	31 Oct 2025 09:42		AR\AJ	Ok
3	PEM	PEM	PL097700.D	31 Oct 2025 09:56		AR\AJ	Ok,M
4	PSTDCCC050	PSTDCCC050	PL097701.D	31 Oct 2025 11:02		AR\AJ	Ok,M
5	Q3488-02	MOO-25-0299	PL097702.D	31 Oct 2025 12:03	F Flag in DCB in both column	AR\AJ	Not Ok
6	Q3490-01	LAW-25-0151	PL097703.D	31 Oct 2025 12:17	TCMX low in 1st column	AR\AJ	Ok,M
7	PB170340BL	PB170340BL	PL097704.D	31 Oct 2025 13:06		AR\AJ	Ok
8	PB170340BS	PB170340BS	PL097705.D	31 Oct 2025 13:19		AR\AJ	Ok
9	Q3495-02	SB-1025-2	PL097706.D	31 Oct 2025 13:33		AR\AJ	Ok,M
10	Q3495-02MS	SB-1025-2MS	PL097707.D	31 Oct 2025 13:47		AR\AJ	Ok,M
11	Q3495-02MSD	SB-1025-2MSD	PL097708.D	31 Oct 2025 14:00		AR\AJ	Ok,M
12	Q3496-01	DRILL-CUTTINGS	PL097709.D	31 Oct 2025 14:14		AR\AJ	Ok,M
13	Q3497-01	OK-01-103025	PL097710.D	31 Oct 2025 14:27		AR\AJ	Ok,M
14	Q3500-01	459	PL097711.D	31 Oct 2025 14:41	DCB high in 1st column	AR\AJ	Ok,M
15	Q3501-01	461	PL097712.D	31 Oct 2025 14:55		AR\AJ	Ok,M
16	Q3502-01	CATCH-CAN-USED-OI	PL097713.D	31 Oct 2025 15:08	DCB high in 1st column	AR\AJ	Ok,M
17	I.BLK	I.BLK	PL097714.D	31 Oct 2025 15:49		AR\AJ	Ok
18	PSTDCCC050	PSTDCCC050	PL097715.D	31 Oct 2025 16:03		AR\AJ	Ok,M

Instrument ID: ECD_L

Daily Analysis Runlog For Sequence/QC Batch ID # PL103125

Review By	rahul	Review On	10/31/2025 2:27:02 PM
Supervise By	yogesh	Supervise On	11/3/2025 8:43:32 AM
SubDirectory	PL103125	HP Acquire Method	HP Processing Method pl102825 8081
STD. NAME	STD REF.#		
Tune/Reschk	PP24942,PP24651		
Initial Calibration Stds	PP24744,PP24746,PP24748,PP24750,PP24751,PP24752,PP24753,PP24755,PP24756,PP24757,PP24758,PP24759,PP24760,PP24761,PP24762,P24763,PP24764		
CCC	PP24751,PP24756,PP24761		
Internal Standard/PEM			
ICV/I.BLK	PP24754,PP24759,PP24761		
Surrogate Standard			
MS/MSD Standard			
LCS Standard			

M : Manual Integration

A
B
C
D
E
F
G
H
I
J
K
L

SOP ID: M3541-ASE Extraction-15

Clean Up SOP #: Florisil

Matrix : Solid

Weigh By: EH

Balance check: EH

Balance ID: EX-SC-2

pH Strip Lot#: N/A

Extraction By: RJ

Filter By: RJ

pH Meter ID: N/A

Hood ID: 3,7

Extraction Start Date : 10/31/2025

Extraction Start Time : 08:28

Extraction End Date : 10/31/2025

Extraction End Time : 11:40

Concentration By: EH

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	PP24766
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	EP2643
Baked Na2SO4	N/A	EP2655
Sand	N/A	E3951
Hexane	N/A	E3981
Florisil	N/A	E3976
9:1 Hexane:Acetone Mixture	N/A	EP2644
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 # 03-40BTS721. Q3500-01, Q3501-01 & Q3502-01 used Limited volume as samples are oil & Oily Debris.

KD Bath ID: N/A

Envap ID: N-EVAP-02

KD Bath Temperature: N/A

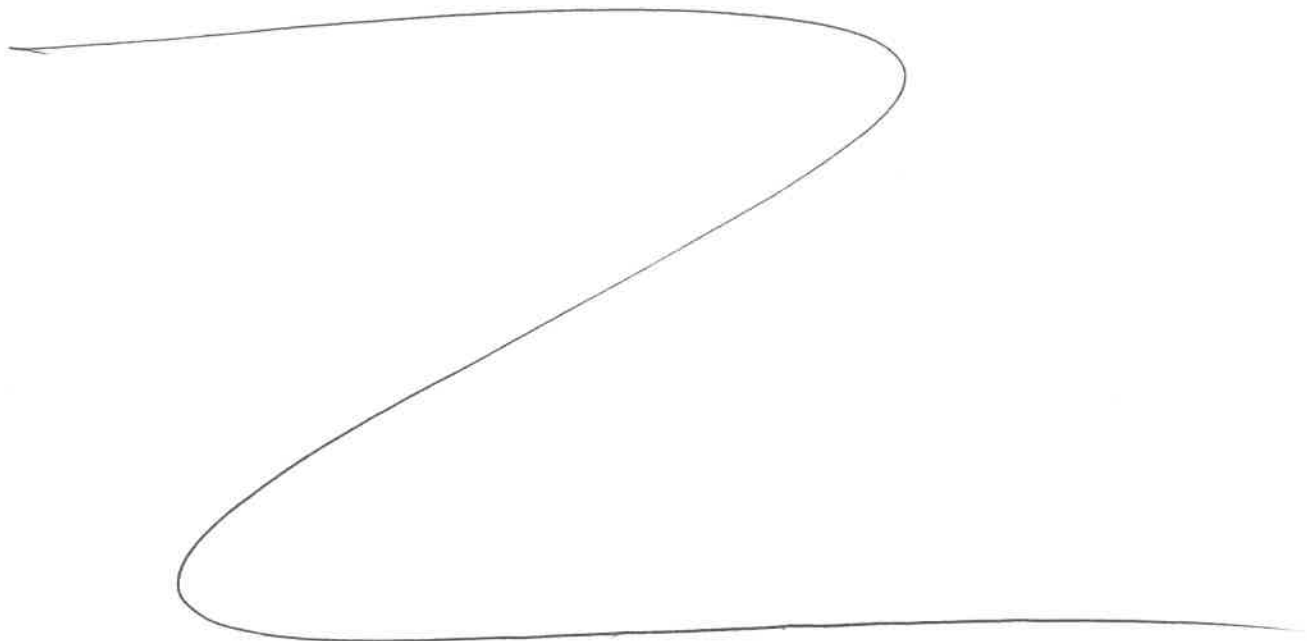
Envap Temperature: 40 °C

Date / Time	Prepped Sample Relinquished By/Location	Received By/Location
10/21/25	RS (Ext Lab)	R.P. Pest PCB
11/25	Preparation Group	Analysis Group

Analytical Method: M3541-ASE Extraction-15

Concentration Date: 10/31/2025

Sample ID	Client Sample ID	Test	g / mL	PH	Surr/Spike By:		Final Vol. (mL)	JarID	Comments	Prep Pos
					AddedBy	VerifiedBy				
PB170340BL	PBLK340	Pesticide-TCL	30.02	N/A	ritesh	Evelyn	10			U1-1
PB170340BS	PLCS340	Pesticide-TCL	30.03	N/A	ritesh	Evelyn	10			2
Q3495-02	SB-1025-2	Pesticide-TCL	30.06	N/A	ritesh	Evelyn	10	E		3
Q3495-02MS	SB-1025-2MS	Pesticide-TCL	30.04	N/A	ritesh	Evelyn	10	E		4
Q3495-02MS D	SB-1025-2MSD	Pesticide-TCL	30.02	N/A	ritesh	Evelyn	10	E		5
Q3496-01	DRILL-CUTTINGS	Pesticide-TCL	30.04	N/A	ritesh	Evelyn	10	E		6
Q3497-01	OK-01-103025	Pesticide-TCL	30.05	N/A	ritesh	Evelyn	10	E		U2-1
Q3500-01	459	Pesticide-TCL	1.06	N/A	ritesh	Evelyn	10		Oil	
Q3501-01	461	Pesticide-TCL	5.03	N/A	ritesh	Evelyn	10	E	Oily Debris	2
Q3502-01	CATCH-CAN-USED-OIL	Pesticide-TCL	1.05	N/A	ritesh	Evelyn	10		Oil	



RS
10/31

17-03-2025 8:23

WORKLIST(Hardcopy Internal Chain)

WorkList Name : Q3495P WorkList ID : 192810 Department : Extraction Date : 10-31-2025 08:23:29

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q3495-02	SB-1025-2	Solid	Pesticide-TCL	Cool 4 deg C	REMI02	D51	10/27/2025	8081B
Q3496-01	DRILL-CUTTINGS	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	D41	10/30/2025	8081B
Q3497-01	OK-01-103025	Solid	Pesticide-TCL	Cool 4 deg C	PSEG05	D41	10/30/2025	8081B
Q3500-01	459	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	D41	10/30/2025	8081B
Q3501-01	461	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	D41	10/30/2025	8081B
Q3502-01	CATCH-CAN-USED-OIL	Solid	Pesticide-TCL	Cool 4 deg C	PSEG03	D41	10/30/2025	8081B

Date/Time 10/31/25 8:23
Raw Sample Received by: R3 CLEA-1006
Raw Sample Relinquished by: JH WDC1

Date/Time 10/31/25 8:50
Raw Sample Received by: JH WDC1
Raw Sample Relinquished by: R3 CLEA-1006



SOP ID: M3510C,3580A-Extraction Pesticide-17

Clean Up SOP #: Florisil **Extraction Start Date :** 10/31/2025

Matrix : Water **Extraction Start Time :** 09:24

Weigh By: N/A **Extraction By:** RS **Extraction End Date :** 10/31/2025

Balance check: N/A **Filter By:** RS **Extraction End Time :** 13:55

Balance ID: N/A **pH Meter ID:** N/A **Concentration By:** EH

pH Strip Lot#: E3953 **Hood ID:** 4,6,7 **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	PP24766
Surrogate	1.0ML	200 PPB	SP24663
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	E3980
Baked Na2SO4	N/A	EP2655
Hexane	N/A	E3981
Florisil	N/A	E3976
9:1 Hexane:Acetone Mixture	N/A	EP2644
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:

40ML Vial Lot#03-40BTS721.

KD Bath ID: WATER BATH-1 **Envap ID:** NEVAP-02

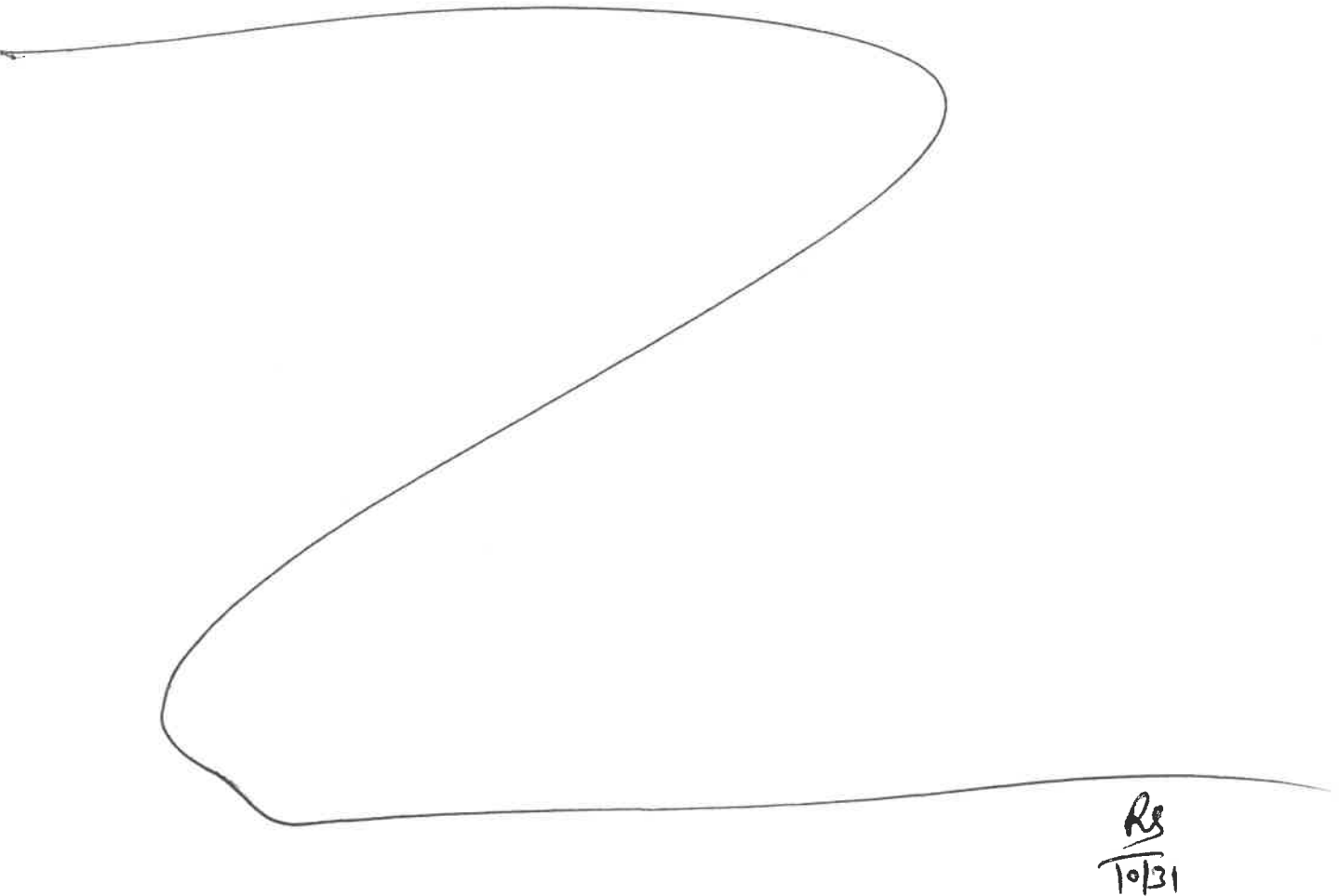
KD Bath Temperature: 60 °C **Envap Temperature:** 40 °C

Date / Time	Prepped Sample Relinquished By/Location	Received By/Location
10/31/25 14:00	RS (Ext Lab)	Y-P-Pesticides
	Preparation Group	Analysis Group

Analytical Method: M3510C,3580A-Extraction Pesticide-17

Concentration Date: 10/31/2025

Sample ID	Client Sample ID	Test	g g mL	PH	Surr/Spike By:		Final Vol. (mL)	JarID	Comments	Prep Pos
					AddedBy	VerifiedBy				
PB170348BL	PBLK348	Pesticide-TCL	1000	6	RUPESH	ritesh	10			SEP-1
PB170348BS	PLCS348	Pesticide-TCL	1000	6	RUPESH	ritesh	10			2
PB170348BS D	PLCSD348	Pesticide-TCL	1000	6	RUPESH	ritesh	10			3
Q3477-03	SOUTH-MAHWAH-WATER	Pesticide-TCL	990	6	RUPESH	ritesh	10	S		4
Q3495-03	TW-1025-1	Pesticide-TCL	980	6	RUPESH	ritesh	10	H		5



Rs
10/31

WORKLIST(Hardcopy Internal Chain)

WorkList Name : Q3477P WorkList ID : 192822 Department : Extraction Date : 10-31-2025 08:43:04

Sample	Customer Sample	Matrix	Test	Preservative	Customer	Raw Sample Storage Location	Collect Date	Method
Q3477-03	SOUTH-MAHWAH-WATER	Water	Pesticide-TCL	Cool 4 deg C	PSEG03	D41	10/28/2025	8081B
Q3495-03	TW-1025-1	Water	Pesticide-TCL	Cool 4 deg C	REMI02	D51	10/27/2025	8081B

Date/Time 10/31/25 9:20
Raw Sample Received by: RS (Ext-Lab)
Raw Sample Relinquished by: RS (Ext-Lab)

Date/Time 10/31/25 9:45
Raw Sample Received by: RS (Ext-Lab)
Raw Sample Relinquished by: RS (Ext-Lab)

LAB CHRONICLE

OrderID:	Q3495	OrderDate:	10/30/2025 10:45:41 AM
Client:	Remington & Vernick Engineers	Project:	Maclearie Park
Contact:	Kyle Carlson	Location:	D51,VOA Ref. #2 Soil,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q3495-02	SB-1025-2	SOIL			10/27/25			10/29/25
			PCB	8082A		10/31/25	10/31/25	
			Pesticide-TCL	8081B		10/31/25	10/31/25	
			EPH_NF	NJEPH		10/31/25	10/31/25	
Q3495-03	TW-1025-1	WATER			10/27/25			10/29/25
			PCB	8082A		10/31/25	10/31/25	
			Pesticide-TCL	8081B		10/31/25	10/31/25	