



SDG NARRATIVE

LAB NAME: Alliance Technical Group, LLC

CASE: 51802 SDG: C0CF7

CONTRACT: 68HERH20D0011

LAB CODE: ACE

LAB ORDER ID: P4627

MODIFICATION REF. NUMBER: NA

Sample ID	EPA Sample ID	pН
P4627-01	C0CN6	
P4627-02	C0CP0	
P4627-03	C0CG2	
P4627-04	C0CR2	
P4627-05	C0CG0	
P4627-06	C0CQ0	
P4627-07	C0CF7	
P4627-08	C0CF2	
P4627-09	C0CF7	
P4627-10	C0CD3	
P4627-11	C0CD9	
P4627-12	C0CD1	
P4627-13	C0CF3	
P4627-14	C0CD2	
P4627-15MS	C0CD2MS	
P4627-16MSD	C0CD2MSD	
P4627-17	C0CF9	
P4627-18	C0CF6	
P4627-19	C0CE0	
P4627-20	C0CF8	
P4627-21	C0CR0	

07 Water samples were delivered to the laboratory intact on 10/30/2024.

Test requested on the Chain of Custody was Semivolatile Organic, Pesticide and Aroclor by Method SFAM01.1.

The temperature of the samples was measured using an I R Gun. The samples temperature was 3.0, 3.6, 3.5, 2.9, 3.3 degree Celsius for the samples received on 10/30/2024. The samples temperature was 2.0, 2.4, 1.7, 1.9, 1.8, 2.3, 2.2, 2.7, 2.5 degree Celsius for the samples received on 10/31/2024.

¹⁴ Water samples were delivered to the laboratory intact on 10/31/2024.





Semivolatiles:

The samples were analyzed on instrument BNA_G using GC Column ZB-GR Semi Volatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGA.

The samples were analyzed on instrument BNA_M using GC Column ZB-GR Semi Volatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGA.

The samples were analyzed on instrument BNA_P using GC Column ZB-GR Semi Volatiles Guardian which is 30 meters, 0.25 mm ID, 0.5 um df, Catalog # 7HG-G027-17-GGA.

Semis volatile Organic sample for water sample was extracted by Method SFAM01.1 on 10/30/2024,10/31/2024, 11/01/2024, The analysis of SVOC-SFAM was based on method SFAM01.1_SVOC.

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria except for,

C0CF8 [4-Nitrophenol-d4 - 9%]. As per method four surrogates are allowed to fail. Therefore no further corrective action was taken.

The Internal Standards Areas met the acceptable requirements.

The Retention Times were acceptable for all samples.

The Blank Spike for {PB164567BS} recoveries met the requirements for all compounds.

The Blank Spike for {PB164571BS} recoveries met the requirements for all compounds.

The Blank Spike for {PB164573BS} recoveries met the requirements for all compounds.

The Blank Spike for {PB164612BS} recoveries met the requirements for all compounds.

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

The Tuning criteria met the requirements.

The Initial Calibration met the requirements.

The Continuous Calibration (SSTD020692) with File ID BP022768.D met the requirements except for Hexachlorocyclopentadiene (-41.2%), under this Continuous Calibration none of the samples were analyzed, therefore no further corrective action was required.

Concentration of Water Sample:

Concentration ug/L = (Ax) (Is) (Vt) (DF) (GPC)

Where.

Ax = Area of the characteristic ion for the compound to be measured.

Ais = Area of the characteristic ion for the internal standard.

Is = Amount of internal standard injected in ng.

Vo = Volume of water extracted in mL.

Vi = Volume of extract injected in uL.

Vt = Volume of the concentrated extract in uL

RRF = Mean Relative Response Factor determined from the initial calibration standard.





GPC = <u>Vin</u> = GPC factor (If no GPC is performed, GPC=1) Vout = Volume of extract collected after GPC cleanup.

Example calculation of C0CF8 for Phenol:

Ax = 25391 Ais = 197858 Is = 20 DF = 1 Vo = 1000 Vi = 1 Vt = 1000 RRF = 1.597 GPC = 1

Concentration ug/L = (25391)(20)(1000)(1)(1)(197858) (1.597) (1000) (1)

= 1.6 ug/L

RRF Calculation of standard 20 ppb for **Naphthalene** with P instrument for method 10/07/2024.

RRF= Area of compound / X Conc. of Internal Standard / Area of Internal Standard Conc. of Compound

= 326983/315808 X 20/20

= 1.035 (Reported RRF)

Pesticides:

The analyses for Pesticides were performed on instrument ECD_D. The front column is ZB-Multi-Residue-2 which is 30 meters, 0.32 mm ID, 0.2 um df. The rear column ZB-Multi-Residue-1 which is 30 meters, 0.32 mm ID, 0.50 um df.

The sample was analyzed on a single injection dual column system. To distinguish the second column analysis from the first column a -2 suffix was added to the file id on the form 1. These refer to forms were both columns are reported. Form 1s for the IBLK and PLCS are referenced as IBLK(1)/IBLK(2), MS(1)/MS(2), MSD(1)/MSD(2) and PLCS01(1) / PLCS01(2) respectively.

Pesticide sample was extracted by method SFAM01.1 on 11/01/2024 and analyzed on 11/01/2024. The sample was extracted and analyzed within contractual holding time.

The Surrogate recoveries met the acceptable criteria.

C0CD2MS met the requirements. C0CD2MSD met the requirements. The RPD met the requirements





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

Blank and Laboratory Control Sample met the requirements.

Retention Times met the requirements.

Florisil check met the requirements.

Resolution Check met the requirements.

The Retention Times were acceptable for all samples.

The Initial Calibration met the requirements.

The Individual Mix A met the requirements.

The Individual Mix B met the requirements.

The PEM met the requirement.

See Manual Integration report for the manual integration information at the end of the case narrative.

Calculation for the Concentration in Water Samples

Concentration ug/L =
$$(Ax) (Vt) (DF) (GPC)$$

(CF) (Vo) (Vi)

Where.

Ax = Response (peak area or height) of the compound to be measured.

CF = Mean Calibration Factor from the initial calibration (area/ng).

Vo = Volume of water extracted in mL.

Vi = Volume of extract injected in uL.

Vt = Volume of the concentrated extract in uL

$$GPC = \underline{Vin} = GPC \text{ factor (If no GPC is performed, GPC=1)}$$

Vout

Vin = Volume of extract loaded onto GPC column.

Vout = Volume of extract collected after GPC cleanup.

Example of Endosulfan I calculation

Calibration Factor Calculation Endosulfan I in the first column

Calibration factor (CF) =
$$\frac{\text{peak area}}{\text{Mass injected in ng}}$$

= $\frac{9516108}{5\text{ng}}$

Mean Calibration Factor = average of 5 point calibration factor

= 1903220

=2025530

No target **Pesticides** were detected in the samples.





Aroclors:

The analyses were performed on instrument GCECD_R. The front column is ZB-MR1 which is 30 meters, 0.32 mm ID, 0.5 um df, Catalogue # 7HM-G016-17. The rear column is ZB-MR2 which is 30 meters, 0.32 mm ID, 0.25 μ m; Catalogue # 7HM-G017-11.

The sample was analyzed on a single injection dual column system. To distinguish the second column analysis from the first column a -2 suffix was added to the file id on the form 1. These refer to forms were both columns are reported. Form 1s for the IBLK and ALCS are referenced as IBLK(1)/IBLK(2), MS(1)/MS(2), MSD(1)/MSD(2) and ALCSO1(1)/ALCSO1(2) respectively.

Aroclor sample was extracted by Method SFAM01.1 on 01/11/2024 and analyzed on 01/11/2024. All the samples were subjected to a Sulfuric acid cleanup. The sample was extracted and analyzed within contractual holding time.

The Surrogate recoveries met the acceptable criteria.

C0CD2MS met the requirements.

C0CD2MSD met the requirements.

The RPD met the requirements

The Laboratory Control Sample met requirements.

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

The Initial Calibration met the requirements.

The Continuing Calibrations met the requirements.

The Retention Times were acceptable for all samples.

See **Manual Integration report** for the manual integration information at the end of the Case narrative.

Calculation for Concentration in Water Samples:

Concentration ug/L =
$$(\underline{Ax})$$
 (Vt) (DF) (GPC)
(CF) (Vo) (Vi)

Where

Ax = Response (peak area or height) of the compound to be measured.

CF = Mean Calibration Factor from the initial calibration (area/ng).

Vo = Volume of water extracted in mL.

Vi = Volume of extract injected in uL.

Vt = Volume of the concentrated extract in uL

GPC = Vin = GPC factor (If no GPC is performed, GPC=1)
Vout

Vin = Volume of extract loaded onto GPC column.

Vout = Volume of extract collected after GPC cleanup.

DF = Dilution Factor.

Example of AR1260 calculation for Peak 1

Calibration factor Peak 1 100ppb ISTD= <u>peak area</u>
Column1 <u>peak area</u>
Mass injected ng





 $= \frac{4574028}{0.100}$

= 45740280 calibration factor for Peak 1 100ppb

Average of 5 peaks = 39861589

No target **Aroclors** were detected in the samples.

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

Signature	Name: Nimisha Pandya.
Date:	Title: Document Control Officer