

## **Report of Analysis**

Client:	Alliance Te	echnical (	Group, LLC - Newa	rk		Date Collected:		
Project:	NJ Waste V	Water PT				Date Received:		
Client Sample I	D: PB168285	BSD				SDG No.:	Q2181	
Lab Sample ID:	: PB1682851	BSD				Matrix:	Water	
Analytical Meth	hod: 8270E					% Solid:	0	
Sample Wt/Vol:		Units:	mL			Final Vol:	1000	uL
Soil Aliquot Vo		emis.	uL			Test:	SVOCMS	
-								S Gloup2
Extraction Type	2:		Decan	ted : N		Level :	LOW	
Injection Volum	ne :		GPC Factor :	1.0		GPC Cleanup :	Ν	PH :
Prep Method :	SW3510C							
File ID/Qc Batch: Dilution:		Prep Date		Date Analyzed		Prep Batch ID		
BF142728.D 1		06/04/25 11:45		06/11/25 11:50		PB168285		
CAS Number	Parameter		Conc.	Qualifier	MDL		LOQ / CRQL	Units
	Parameter		Conc.	Qualifier	MDL		LOQ / CRQL	Units
CAS Number TARGETS 105-67-9	Parameter 2,4-Dimethylphenol		<b>Conc.</b> 46.3	Qualifier	<b>MDL</b> 1.90		LOQ / CRQL 5.00	Units ug/L
<b>TARGETS</b> 105-67-9				Qualifier				
TARGETS				Qualifier				
TARGETS 105-67-9 SURROGATES	2,4-Dimethylphenol		46.3	Qualifier	1.90		5.00	ug/L
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4	2,4-Dimethylphenol 2-Fluorophenol	ol	46.3 125	Qualifier	1.90 23 - 138		5.00 83%	ug/L SPK: 150
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4 13127-88-3	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno	ol	46.3 125 127	Qualifier	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4 13127-88-3 118-79-6	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno		46.3 125 127	<b>Qualifier</b> 6.893	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4 13127-88-3 118-79-6 <b>INTERNAL STAN</b>	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno NDARDS		46.3 125 127 129	-	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4 13127-88-3 118-79-6 <b>INTERNAL STAN</b> 3855-82-1	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno NDARDS 1,4-Dichlorobenzene		46.3 125 127 129 80000	6.893	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150
TARGETS   105-67-9   SURROGATES   367-12-4   13127-88-3   118-79-6   INTERNAL STAN   3855-82-1   1146-65-2	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno <b>NDARDS</b> 1,4-Dichlorobenzene Naphthalene-d8		46.3 125 127 129 80000 305000	6.893 8.181	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150
<b>TARGETS</b> 105-67-9 <b>SURROGATES</b> 367-12-4 13127-88-3 118-79-6 <b>INTERNAL STAN</b> 3855-82-1 1146-65-2 15067-26-2	2,4-Dimethylphenol 2-Fluorophenol Phenol-d6 2,4,6-Tribromopheno NDARDS 1,4-Dichlorobenzene Naphthalene-d8 Acenaphthene-d10		46.3 125 127 129 80000 305000 168000	6.893 8.181 9.939	1.90 23 - 138 10 - 134		5.00 83% 84%	ug/L SPK: 150 SPK: 150

U = Not Detected

- LOQ = Limit of Quantitation
- MDL = Method Detection Limit
- LOD = Limit of Detection
- E = Value Exceeds Calibration Range
- 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
- () = Laboratory InHouse Limit
- A = Aldol-Condensation Reaction Products