

Report of Analysis

Client: PSEG
Project: Orange Gas
Client Sample ID: SB-18DL
Lab Sample ID: Q3446-05DL
Analytical Method: 8260D
Sample Wt/Vol: 5 mL

Level: LOW
Final Vol: 5000 uL

Date Collected: 10/23/25
Date Received: 10/23/25
SDG No.: Q3446
Matrix: TCLP
% Solid: 0
Test: TCLP VOA

CAS Number	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Date Ana.	BatchID
TARGETS									
75-01-4	Vinyl Chloride	0.25	UD	50	0.013	0.25	mg/L	10/27/25 15:23	VX102725
75-35-4	1,1-Dichloroethene	0.25	UD	50	0.012	0.25	mg/L	10/27/25 15:23	VX102725
78-93-3	2-Butanone	1.30	UD	50	0.049	1.30	mg/L	10/27/25 15:23	VX102725
56-23-5	Carbon Tetrachloride	0.25	UD	50	0.013	0.25	mg/L	10/27/25 15:23	VX102725
67-66-3	Chloroform	0.25	UD	50	0.013	0.25	mg/L	10/27/25 15:23	VX102725
71-43-2	Benzene	1.40	D	50	0.0075	0.25	mg/L	10/27/25 15:23	VX102725
107-06-2	1,2-Dichloroethane	0.25	UD	50	0.011	0.25	mg/L	10/27/25 15:23	VX102725
79-01-6	Trichloroethene	0.25	UD	50	0.0047	0.25	mg/L	10/27/25 15:23	VX102725
127-18-4	Tetrachloroethene	0.25	UD	50	0.012	0.25	mg/L	10/27/25 15:23	VX102725
108-90-7	Chlorobenzene	0.25	UD	50	0.0060	0.25	mg/L	10/27/25 15:23	VX102725
SURROGATES									
17060-07-0	1,2-Dichloroethane-d4	56.1			74 - 125	112%	SPK: 50		
1868-53-7	Dibromofluoromethane	40.5			75 - 124	81%	SPK: 50		
2037-26-5	Toluene-d8	47.4			86 - 113	95%	SPK: 50		
460-00-4	4-Bromofluorobenzene	52.8			77 - 121	106%	SPK: 50		
INTERNAL STANDARDS									
		Area Count							
363-72-4	Pentafluorobenzene	154000							
540-36-3	1,4-Difluorobenzene	309000							
3114-55-4	Chlorobenzene-d5	307000							
3855-82-1	1,4-Dichlorobenzene-d4	152000							

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