

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062425\
 Data File : VU063434.D
 Acq On : 24 Jun 2025 12:54
 Operator : MD/SY
 Sample : Q2377-01
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 PW-B6-L66-061925

Quant Time: Jun 25 08:21:44 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U062325DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Jun 25 03:53:43 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Fluorobenzene	6.107	96	46278	1.000	ug/l	# 0.00
System Monitoring Compounds						
57) 4-Bromofluorobenzene	10.634	95	13619	0.802	ug/l	0.00
Spiked Amount 1.000			Recovery	=	80.000%	
68) 1,2-Dichlorobenzene-d4	12.190	152	13641	0.799	ug/l	0.00
Spiked Amount 1.000			Recovery	=	80.000%	
Target Compounds						
						Qvalue
3) Chloromethane	1.522	50	4919	0.210	ug/l	99
27) Chloroform	5.071	83	1047169	38.909	ug/l	99
44) Bromodichloromethane	7.097	83	223947	11.817	ug/l	99
55) Dibromochloromethane	8.804	129	25656	2.186	ug/l	98
64) m/p-Xylenes	9.695	106	618	0.809	ug/l #	64
78) 1,2,4-Trimethylbenzene	11.473	105	1007	0.434	ug/l	98
81) p-Isopropyltoluene	11.791	119	1272	0.432	ug/l	92
89) Naphthalene	14.094	128	9003	1.136	ug/l #	93

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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