

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030625\
 Data File : VX045176.D
 Acq On : 06 Mar 2025 18:12
 Operator : JC/MD
 Sample : Q1490-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CLAY-SLUDGE-DRUMS

Quant Time: Mar 07 00:48:58 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
 Quant Title : SW846 8260
 QLast Update : Fri Feb 28 06:45:16 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	79920	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	159351	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	146313	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	59972	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	67372	52.997	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 106.000%		
35) Dibromofluoromethane	5.379	113	54432	51.084	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.160%		
50) Toluene-d8	8.647	98	197597	51.152	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 102.300%		
62) 4-Bromofluorobenzene	11.079	95	64828	50.644	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 101.280%		
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	57468	97.433	ug/l	99
18) Methyl Acetate	2.709	43	3719	2.621	ug/l	94
20) Methylene Chloride	2.788	84	2376	2.034	ug/l	94
25) 2-Butanone	4.574	43	7989	8.998	ug/l	93
43) Isopropyl Acetate	6.342	43	45459	16.327	ug/l	98
52) Toluene	8.720	92	6073	2.243	ug/l	96

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

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