

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
 Data File : VX043999.D
 Acq On : 25 Nov 2024 20:03
 Operator : JC/MD
 Sample : P4938-04
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-732

Quant Time: Nov 26 00:55:33 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 22 03:45:52 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	117541	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	225495	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	193570	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	87608	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	100223	49.713	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	99.420%		
35) Dibromofluoromethane	5.379	113	71907	44.627	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	89.260%		
50) Toluene-d8	8.647	98	269999	48.611	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	97.220%		
62) 4-Bromofluorobenzene	11.079	95	92663	49.021	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	98.040%		
Target Compounds						
16) Acetone	2.392	43	42513	45.913	ug/l	97
18) Methyl Acetate	2.709	43	3560	1.651	ug/l #	87
20) Methylene Chloride	2.788	84	2244	1.428	ug/l #	93
43) Isopropyl Acetate	6.342	43	115619	29.149	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112524\
Data File : VX043999.D
Acq On : 25 Nov 2024 20:03
Operator : JC/MD
Sample : P4938-04
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 24 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-732

Quant Time: Nov 26 00:55:33 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X112124W.M
Quant Title : SW846 8260
QLast Update : Fri Nov 22 03:45:52 2024
Response via : Initial Calibration

