

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112624\
 Data File : VX044015.D
 Acq On : 26 Nov 2024 14:27
 Operator : JC/MD
 Sample : P5003-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW-10A-HYD-20241125

Quant Time: Nov 27 00:53:36 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	5.544	168	106926	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	202031	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	179566	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	76328	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	92084	50.210	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	100.420%
35) Dibromofluoromethane	5.385	113	68197	47.240	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.480%
50) Toluene-d8	8.647	98	249740	50.186	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.380%
62) 4-Bromofluorobenzene	11.079	95	83026	49.024	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	98.040%
Target Compounds						
					Qvalue	
3) Chloromethane	1.294	50	444	0.311	ug/l #	82
16) Acetone	2.380	43	1596	1.895	ug/l	92
60) Dibromochloromethane	9.525	129	1062	0.765	ug/l	87

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112624\
Data File : VX044015.D
Acq On : 26 Nov 2024 14:27
Operator : JC/MD
Sample : P5003-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
RW-10A-HYD-20241125

Quant Time: Nov 27 00:53:36 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

