

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX123124\
 Data File : VX044555.D
 Acq On : 30 Dec 2024 19:38
 Operator : JC/MD
 Sample : P5376-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RW10A-MW01S-20241218

Quant Time: Dec 31 00:36:51 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121124W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 12 04:28:14 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	122970	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	208916	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	183505	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	75781	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	91534	42.454	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	84.900%		
35) Dibromofluoromethane	5.385	113	72294	49.964	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	99.920%		
50) Toluene-d8	8.647	98	261648	53.599	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	107.200%		
62) 4-Bromofluorobenzene	11.079	95	84375	49.466	ug/l	0.00
Spiked Amount 50.000	Range 77 - 121		Recovery =	98.940%		
Target Compounds						
3) Chloromethane	1.300	50	1022	0.562	ug/l	97
16) Acetone	2.392	43	1176	1.338	ug/l #	81
20) Methylene Chloride	2.782	84	559	0.339	ug/l #	89
22) Diisopropyl ether	3.776	45	9268	1.793	ug/l #	80

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

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