

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121024\
 Data File : VX044215.D
 Acq On : 10 Dec 2024 17:57
 Operator : JC/MD
 Sample : P5246-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW7-20241209

Quant Time: Dec 11 01:39:37 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	5.549	168	107458	50.000	ug/l	0.01
34) 1,4-Difluorobenzene	6.757	114	207402	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	185141	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	78742	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	97507	52.904	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	105.800%
35) Dibromofluoromethane	5.385	113	69956	47.204	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.400%
50) Toluene-d8	8.646	98	254455	49.809	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.620%
62) 4-Bromofluorobenzene	11.079	95	86650	49.839	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	99.680%
Target Compounds						
					Qvalue	
3) Chloromethane	1.300	50	493	0.344	ug/l #	87
16) Acetone	2.391	43	967	1.142	ug/l #	83
20) Methylene Chloride	2.794	84	469	0.326	ug/l #	75
30) Chloroform	5.092	83	13000	4.843	ug/l	94

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

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