

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101824\
 Data File : VX043442.D
 Acq On : 18 Oct 2024 18:14
 Operator : JC/MD
 Sample : P4440-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BPOW6-11-HYD-20241016

Quant Time: Oct 18 23:24:44 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X100124W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 02 16:50:57 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	45096	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	82530	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	68771	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	27812	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	34503	46.848	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	93.700%
35) Dibromofluoromethane	5.385	113	26061	45.789	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.580%
50) Toluene-d8	8.647	98	96994	49.176	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.360%
62) 4-Bromofluorobenzene	11.079	95	33026	46.090	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	92.180%
Target Compounds						
					Qvalue	
30) Chloroform	5.093	83	1239	1.170	ug/l #	74
47) Bromodichloromethane	7.824	83	1924	2.283	ug/l #	94
60) Dibromochloromethane	9.525	129	1880	2.984	ug/l	90
71) Bromoform	10.805	173	609	1.521	ug/l #	91

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

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