

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX100924\
 Data File : VX043343.D
 Acq On : 09 Oct 2024 16:40
 Operator : JC/MD
 Sample : P4365-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BPOW6-11-HYD-20241008

Quant Time: Oct 10 01:34:24 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	5.550	168	83434	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	156513	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	133681	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	55217	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	72840	53.456	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.920%	
35) Dibromofluoromethane	5.385	113	52265	48.422	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 96.840%	
50) Toluene-d8	8.647	98	186223	49.785	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 99.580%	
62) 4-Bromofluorobenzene	11.079	95	64380	47.376	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 94.760%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	1417	2.490	ug/l #	88
18) Methyl Acetate	2.709	43	1819	1.274	ug/l	95
47) Bromodichloromethane	7.824	83	2597	1.625	ug/l #	97
60) Dibromochloromethane	9.525	129	2869	2.402	ug/l	94
71) Bromoform	10.799	173	1076	1.382	ug/l #	94

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

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