

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080924\
 Data File : VX042998.D
 Acq On : 09 Aug 2024 15:50
 Operator : JC/MD
 Sample : P3515-03RE
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BPOW6-7-HYD-20240806RE

Quant Time: Aug 09 23:27:04 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X080724W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 07 02:55:22 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	154029	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	246282	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	224835	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	97197	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109590	51.828	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 103.660%		
35) Dibromofluoromethane	5.391	113	83006	51.547	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 103.100%		
50) Toluene-d8	8.647	98	304018	53.181	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 106.360%		
62) 4-Bromofluorobenzene	11.079	95	98964	48.175	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 96.360%		
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	3075	3.353	ug/l #	89
30) Chloroform	5.105	83	3393	1.000	ug/l #	76
47) Bromodichloromethane	7.824	83	5664	2.415	ug/l	89
60) Dibromochloromethane	9.525	129	7986	4.351	ug/l	99
71) Bromoform	10.805	173	5437	4.792	ug/l #	92

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

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