

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051923\
 Data File : VX035767.D
 Acq On : 19 May 2023 17:50
 Operator : JC/MD
 Sample : 02860-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-TB-03-05182023

Quant Time: May 22 05:19:20 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 11 13:57:54 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	183687	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	356947	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	313740	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	135152	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	183798	57.356	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 114.720%	
35) Dibromofluoromethane	5.385	113	112569	46.763	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 93.520%	
50) Toluene-d8	8.647	98	424875	47.777	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 95.560%	
62) 4-Bromofluorobenzene	11.079	95	162218	45.304	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 90.600%	
Target Compounds						
					Qvalue	
16) Acetone	2.404	43	5291	3.842	ug/l	90
29) Tetrahydrofuran	5.038	42	1950	1.627	ug/l	# 88
37) Ethyl Acetate	4.715	43	43944	11.088	ug/l	99
40) Benzene	6.038	78	2947	0.277	ug/l	98
44) Trichloroethene	7.123	130	2054	0.762	ug/l	84
52) Toluene	8.720	92	8573	1.278	ug/l	95
67) Ethyl Benzene	10.195	91	2863	0.222	ug/l	89
68) m/p-Xylenes	10.299	106	4419	0.928	ug/l	94
69) o-Xylene	10.640	106	2321	0.491	ug/l	91
84) 1,2,4-Trimethylbenzene	11.750	105	2527	0.260	ug/l	96

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

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