

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051923\
 Data File : VX035771.D
 Acq On : 19 May 2023 19:23
 Operator : JC/MD
 Sample : 02860-17
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-EQB-12-05172023

Quant Time: May 22 05:20:34 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	178473	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	347371	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	308808	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	134222	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	182743	58.693	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	117.380%#
35) Dibromofluoromethane	5.385	113	115928	49.486	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.980%
50) Toluene-d8	8.647	98	413960	47.832	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.660%
62) 4-Bromofluorobenzene	11.079	95	160493	46.058	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	92.120%
Target Compounds						
						Qvalue
16) Acetone	2.392	43	3203	2.394	ug/l	# 84
29) Tetrahydrofuran	5.019	42	21003	18.039	ug/l	95
30) Chloroform	5.092	83	183630	40.191	ug/l	100
47) Bromodichloromethane	7.817	83	4947	1.365	ug/l	95
52) Toluene	8.720	92	1449	0.222	ug/l	84

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

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