

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX050523\
 Data File : VX035542.D
 Acq On : 05 May 2023 17:35
 Operator : JC/MD
 Sample : 02643-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-IDW-AQ-05042023

Quant Time: May 08 00:52:38 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X042523W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 25 16:31:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	286285	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	474638	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	429349	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	200742	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	172191	49.855	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	99.700%		
35) Dibromofluoromethane	5.391	113	121101	41.570	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	83.140%		
50) Toluene-d8	8.647	98	559755	50.935	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	101.880%		
62) 4-Bromofluorobenzene	11.079	95	209775	50.743	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	101.480%		
Target Compounds					Qvalue	
16) Acetone	2.404	43	693570	462.359	ug/l	98
30) Chloroform	5.099	83	23445	4.258	ug/l	95
44) Trichloroethene	7.129	130	16142	4.490	ug/l	98
60) Dibromochloromethane	9.525	129	4691	1.485	ug/l	98
71) Bromoform	10.805	173	2324	1.034	ug/l #	97

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

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