

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091521\
 Data File : VX024270.D
 Acq On : 15 Sep 2021 18:59
 Operator : JC/MD
 Sample : M3753-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-1

Quant Time: Sep 23 09:00:09 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091421W.M
 Quant Title : SW846 8260
 QLast Update : Tue Sep 21 10:52:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	129226	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	223952	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	181761	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	83294	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	95678	51.922	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery = 103.840%			
35) Dibromofluoromethane	5.397	113	73709	49.077	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.160%			
50) Toluene-d8	8.653	98	273517	49.990	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery = 99.980%			
62) 4-Bromofluorobenzene	11.085	95	83780	39.346	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 78.700%#			
Target Compounds						Qvalue
3) Chloromethane	1.294	50	913	0.704	ug/l	91
6) Chloroethane	1.691	64	2318	2.718	ug/l	98
16) Acetone	2.386	43	95019	120.932	ug/l	92
20) Methylene Chloride	2.788	84	1920	1.273	ug/l #	77
25) 2-Butanone	4.574	43	6194	5.352	ug/l	94
30) Chloroform	5.099	83	935	0.352	ug/l #	75
42) 1,2-Dichloroethane	6.098	62	2141	0.869	ug/l	86
52) Toluene	8.720	92	9910	2.560	ug/l	98
55) 1,1,2-Trichloroethane	9.153	97	934	0.567	ug/l #	85

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

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