

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX121322\
 Data File : VX033391.D
 Acq On : 13 Dec 2022 14:44
 Operator : JC/MD
 Sample : N6001-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-6M

Quant Time: Dec 14 05:38:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 07 12:43:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.555	168	139160	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	221430	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	196492	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	90787	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	41641	46.643	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	93.280%		
35) Dibromofluoromethane	5.385	113	47845	48.040	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.080%		
50) Toluene-d8	8.652	98	98580	47.315	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	94.640%		
62) 4-Bromofluorobenzene	11.079	95	75631	47.475	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	94.960%		
Target Compounds					Qvalue	
16) Acetone	2.391	43	1338	2.025	ug/l	98
30) Chloroform	5.098	83	4242	1.507	ug/l	94
40) Benzene	6.049	78	7675	1.356	ug/l	92
52) Toluene	8.720	92	2204	0.602	ug/l	93
67) Ethyl Benzene	10.195	91	15172	2.189	ug/l	98
68) m/p-Xylenes	10.305	106	13561	5.028	ug/l	97
69) o-Xylene	10.640	106	2709	1.016	ug/l	87
80) 1,3,5-Trimethylbenzene	11.451	105	1239	0.226	ug/l	99
84) 1,2,4-Trimethylbenzene	11.756	105	3625	0.663	ug/l	98

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

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