

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072622\
 Data File : VX030210.D
 Acq On : 26 Jul 2022 15:16
 Operator : JC/MD
 Sample : N3876-01 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 SPC-02-84

Quant Time: Jul 26 17:21:46 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 25 16:49:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	143925	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	245553	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	251281	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	139665	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	151886	52.505	ug/l	-0.01
Spiked Amount	50.000	Range	74 - 125	Recovery	= 105.000%	
35) Dibromofluoromethane	5.385	113	122471	52.174	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.340%	
50) Toluene-d8	8.647	98	459874	50.588	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 101.180%	
62) 4-Bromofluorobenzene	11.085	95	160789	47.659	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.320%	
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.941	59	5935	8.411	ug/l #	77
16) Acetone	2.386	43	2537	1.713	ug/l #	73
17) Carbon Disulfide	2.508	76	1737	0.286	ug/l #	72
30) Chloroform	5.093	83	1079	0.234	ug/l #	53
44) Trichloroethene	7.129	130	890	0.306	ug/l	90
64) Tetrachloroethene	9.275	164	207586	76.735	ug/l	97
90) Hexachloroethane	12.542	117	28089	17.606	ug/l	97

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

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