

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX080723\
 Data File : VX036849.D
 Acq On : 07 Aug 2023 20:05
 Operator : JC/MD
 Sample : 03903-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-01W

Quant Time: Aug 08 01:41:26 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072123W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 21 12:14:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	128225	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	234778	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	221835	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	105195	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	89656	47.255	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	94.520%		
35) Dibromofluoromethane	5.391	113	75989	47.089	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.180%		
50) Toluene-d8	8.646	98	286749	50.108	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.220%		
62) 4-Bromofluorobenzene	11.079	95	115328	53.416	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	106.840%		
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	7639	9.703	ug/l	93
17) Carbon Disulfide	2.513	76	1716	0.433	ug/l	96
31) Cyclohexane	5.495	56	9599	3.654	ug/l #	57
39) Methylcyclohexane	7.378	83	46755	15.633	ug/l #	86
67) Ethyl Benzene	10.195	91	5397	0.604	ug/l	99
73) Isopropylbenzene	10.963	105	71968	8.267	ug/l	99
78) n-propylbenzene	11.305	91	76693	7.617	ug/l	99
83) tert-Butylbenzene	11.713	119	13210	1.846	ug/l	89
84) 1,2,4-Trimethylbenzene	11.750	105	27907	3.831	ug/l	96
85) sec-Butylbenzene	11.890	105	181457	20.336	ug/l	97
89) n-Butylbenzene	12.329	91	119053	18.114	ug/l #	67

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

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