

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072122\
 Data File : VX030149.D
 Acq On : 20 Jul 2022 21:53
 Operator : JC/MD
 Sample : N3760-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 AUD-1821

Quant Time: Jul 21 03:34:37 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X071222W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 13 09:05:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	176306	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	325269	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	282309	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	120326	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	117301	54.661	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.320%	
35) Dibromofluoromethane	5.391	113	96850	46.609	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 93.220%	
50) Toluene-d8	8.653	98	376396	48.163	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 96.320%	
62) 4-Bromofluorobenzene	11.085	95	124131	45.770	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 91.540%	
Target Compounds						
						Qvalue
16) Acetone	2.380	43	214061	211.692	ug/l	93
17) Carbon Disulfide	2.514	76	9193	1.463	ug/l	98
18) Methyl Acetate	2.709	43	1114	0.362	ug/l	91
25) 2-Butanone	4.568	43	54958	31.906	ug/l	94
51) 4-Methyl-2-Pentanone	8.580	43	4967	1.447	ug/l	98
86) p-Isopropyltoluene	12.012	119	3871	0.539	ug/l #	82
95) Naphthalene	13.780	128	6302	0.657	ug/l #	92

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

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