

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040622\
 Data File : VN071892.D
 Acq On : 06 Apr 2022 20:20
 Operator : JC\MD
 Sample : N2175-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 023-OUTFALL

Quant Time: Apr 07 04:31:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	628751	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1082779	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1117357	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	410797	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	340387	47.421	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	94.840%		
35) Dibromofluoromethane	8.022	113	307466	48.256	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.520%		
50) Toluene-d8	10.439	98	1361339	52.230	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	104.460%		
62) 4-Bromofluorobenzene	12.727	95	463835	55.024	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	110.040%		
Target Compounds						
					Qvalue	
13) Acrolein	4.040	56	4541	3.197	ug/l	97
16) Acetone	4.293	43	93006	30.319	ug/l	97
18) Methyl Acetate	4.863	43	15572	0.553	ug/l #	69
25) 2-Butanone	7.339	43	39348	8.898	ug/l	99
51) 4-Methyl-2-Pentanone	10.327	43	14461	1.553	ug/l	98
60) Dibromochloromethane	11.233	129	2983	0.418	ug/l	88
70) Styrene	12.292	104	41396	1.891	ug/l	90

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

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