

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN083122\
 Data File : VN074213.D
 Acq On : 31 Aug 2022 14:45
 Operator : JC\MD
 Sample : N4350-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMMUNITY-CENTER-KITCHEN-SINK

Quant Time: Sep 01 00:25:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 31 07:02:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	214190	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	392477	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	381707	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	169356	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	166185	50.016	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.040%		
35) Dibromofluoromethane	7.951	113	138672	54.298	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 108.600%		
50) Toluene-d8	10.380	98	456759	45.414	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 90.820%		
62) 4-Bromofluorobenzene	12.674	95	181975	59.496	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 119.000%		
Target Compounds						
						Qvalue
3) Chloromethane	2.269	50	1740	0.557	ug/l #	78
29) Tetrahydrofuran	7.610	42	86091	38.061	ug/l #	81
30) Chloroform	7.745	83	14300	2.492	ug/l	92
47) Bromodichloromethane	9.698	83	10207	2.220	ug/l #	96
60) Dibromochloromethane	11.180	129	9937	3.198	ug/l	99
71) Bromoform	12.392	173	4877	1.886	ug/l #	90

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

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