

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074489.D
 Acq On : 16 Sep 2022 03:16
 Operator : JC\MD
 Sample : N4630-17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-17-4-4

Quant Time: Sep 16 07:01:25 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091522W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 19:20:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	389573	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	754570	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	768698	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	337993	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	285022	43.120	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	86.240%		
35) Dibromofluoromethane	7.951	113	240004	47.881	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	95.760%		
50) Toluene-d8	10.380	98	897878	46.318	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	92.640%		
62) 4-Bromofluorobenzene	12.674	95	358742	54.033	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	108.060%		
Target Compounds						Qvalue
16) Acetone	4.216	43	160776	38.011	ug/l	99
20) Methylene Chloride	5.022	84	23354	3.645	ug/l #	85
37) Ethyl Acetate	7.334	43	81739	5.966	ug/l #	96
43) Isopropyl Acetate	8.492	43	367849	17.139	ug/l #	82
84) 1,2,4-Trimethylbenzene	13.310	105	70192	2.472	ug/l	99

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

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