

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074487.D
 Acq On : 16 Sep 2022 02:28
 Operator : JC\MD
 Sample : N4630-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-17-4-2-GR

Quant Time: Sep 16 07:01:01 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	396999	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	768144	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	794202	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	353396	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	285756	42.422	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.840%
35) Dibromofluoromethane	7.951	113	241898	47.406	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.820%
50) Toluene-d8	10.380	98	898553	45.534	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.060%
62) 4-Bromofluorobenzene	12.674	95	370628	54.837	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	109.680%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	101714	23.597	ug/l	94
18) Methyl Acetate	4.787	43	13390	1.000	ug/l #	88
20) Methylene Chloride	5.010	84	20840	3.196	ug/l	95
37) Ethyl Acetate	7.333	43	99447	7.130	ug/l	95
43) Isopropyl Acetate	8.492	43	485311	22.213	ug/l #	82
84) 1,2,4-Trimethylbenzene	13.310	105	38800	1.307	ug/l	93

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

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