

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081122\
 Data File : VN073835.D
 Acq On : 11 Aug 2022 19:21
 Operator : JC\MD
 Sample : N4120-07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BLDG-19C

Quant Time: Aug 12 06:09:13 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 10 05:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	418115	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	663641	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	649505	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	310947	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	241402	47.376	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	94.760%
35) Dibromofluoromethane	7.951	113	227740	54.650	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	109.300%
50) Toluene-d8	10.380	98	748615	45.651	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	91.300%
62) 4-Bromofluorobenzene	12.674	95	283081	49.220	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	98.440%
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	102526	37.462	ug/l	99
18) Methyl Acetate	4.793	43	8020	1.178	ug/l #	87
20) Methylene Chloride	5.016	84	30117	4.810	ug/l	99
25) 2-Butanone	7.269	43	30364	7.144	ug/l	92
43) Isopropyl Acetate	8.498	43	53240	5.039	ug/l #	92

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

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