

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122121\
 Data File : VN070279.D
 Acq On : 21 Dec 2021 20:07
 Operator : JC/MD
 Sample : M5142-02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 A2

Quant Time: Dec 21 21:01:59 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
 Quant Title : SW846 8260
 QLast Update : Fri Dec 03 08:46:47 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	670536	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	1238967	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1059092	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	355720	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	684592	66.124	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 132.240%	
35) Dibromofluoromethane	8.024	113	437419	56.364	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 112.720%	
50) Toluene-d8	10.440	98	1478985	47.462	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 94.920%	
62) 4-Bromofluorobenzene	12.731	95	487024	43.141	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 86.280%	
Target Compounds						
					Qvalue	
16) Acetone	4.291	43	240277	38.581	ug/l	96
18) Methyl Acetate	4.865	43	29498	1.569	ug/l #	69
20) Methylene Chloride	5.093	84	83829	10.125	ug/l	89
25) 2-Butanone	7.340	43	41928	4.955	ug/l #	88
43) Isopropyl Acetate	8.560	43	157999	6.082	ug/l #	84
67) Ethyl Benzene	11.843	91	294352	7.542	ug/l	96
68) m/p-Xylenes	11.948	106	452449	31.839	ug/l	89
69) o-Xylene	12.280	106	169965	11.874	ug/l	90
78) n-propylbenzene	12.919	91	35225	1.000	ug/l	96
80) 1,3,5-Trimethylbenzene	13.058	105	81011	3.180	ug/l	97
84) 1,2,4-Trimethylbenzene	13.366	105	288854	11.907	ug/l	98

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

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