

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
 Data File : VN069576.D
 Acq On : 19 Nov 2021 2:23
 Operator : JC/MD
 Sample : PB140886ZHE#02
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB140886ZHE#02

Quant Time: Nov 19 05:15:34 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N110921W.M
 Quant Title : SW846 8260
 QLast Update : Fri Nov 19 04:51:32 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	1321712	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.968	114	2381557	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2448709	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	934821	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	987251	51.516	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 103.040%		
35) Dibromofluoromethane	8.024	113	679346	47.561	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 95.120%		
50) Toluene-d8	10.443	98	3040624	54.126	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 108.260%		
62) 4-Bromofluorobenzene	12.734	95	1109634	53.768	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 107.540%		
Target Compounds						
					Qvalue	
16) Acetone	4.293	43	225938	21.851	ug/l	94
18) Methyl Acetate	4.865	43	48116	1.432	ug/l #	78
20) Methylene Chloride	5.098	84	41190	2.809	ug/l #	86
25) 2-Butanone	7.351	43	61909	4.422	ug/l	91
29) Tetrahydrofuran	7.702	42	23793m	2.769	ug/l	
43) Isopropyl Acetate	8.566	43	193544	4.448	ug/l #	83

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

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