

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN111821\
 Data File : VN069575.D
 Acq On : 19 Nov 2021 1:58
 Operator : JC/MD
 Sample : PB140886ZHE#01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB140886ZHE#01

Quant Time: Nov 19 05:15:21 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	1313901	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	2350108	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	2387846	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	920412	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	969181	50.873	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 101.740%	
35) Dibromofluoromethane	8.024	113	674961	47.887	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 95.780%	
50) Toluene-d8	10.443	98	2985960	53.864	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 107.720%	
62) 4-Bromofluorobenzene	12.733	95	1073806	52.728	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 105.460%	
Target Compounds						
					Qvalue	
16) Acetone	4.296	43	227884	22.170	ug/l	93
18) Methyl Acetate	4.867	43	37302	1.116	ug/l #	78
20) Methylene Chloride	5.101	84	41514	2.848	ug/l	88
25) 2-Butanone	7.348	43	65369	4.697	ug/l #	90
29) Tetrahydrofuran	7.697	42	24442	2.862	ug/l	88
43) Isopropyl Acetate	8.566	43	181662	4.231	ug/l #	84

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

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