

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN122822\
 Data File : VN076021.D
 Acq On : 29 Dec 2022 00:31
 Operator : JC\MD
 Sample : PB149923ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB149923ZHE#02

Quant Time: Dec 29 07:48:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N121422W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 14 13:03:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	378574	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	597810	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	527459	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	220684	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	190058	46.147	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	92.300%		
35) Dibromofluoromethane	8.171	113	174028	47.007	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.020%		
50) Toluene-d8	10.571	98	696421	50.312	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	100.620%		
62) 4-Bromofluorobenzene	12.853	95	218727	47.709	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	95.420%		
Target Compounds						Qvalue
16) Acetone	4.447	43	67492	47.538	ug/l	97
20) Methylene Chloride	5.289	84	19179	4.467	ug/l	97
25) 2-Butanone	7.494	43	24405	12.277	ug/l	92
37) Ethyl Acetate	7.571	43	36621	9.270	ug/l	98
43) Isopropyl Acetate	8.694	43	286953	44.036	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	22773	5.779	ug/l #	75

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

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