

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078559.D
 Acq On : 18 Jul 2023 14:25
 Operator : JC\MD
 Sample : PB154251ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154251ZHE#04

Quant Time: Jul 19 01:50:00 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	408037	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	748819	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	738854	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	287021	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	274074	50.103	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 100.200%		
35) Dibromofluoromethane	8.177	113	277529	54.597	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.200%		
50) Toluene-d8	10.571	98	923587	48.096	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 96.200%		
62) 4-Bromofluorobenzene	12.853	95	319291	53.981	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 107.960%		
Target Compounds						
						Qvalue
16) Acetone	4.442	43	115077	67.535	ug/l	91
20) Methylene Chloride	5.289	84	27298	4.472	ug/l	93
25) 2-Butanone	7.500	43	83050	30.963	ug/l #	85
37) Ethyl Acetate	7.571	43	22820	3.730	ug/l	99
43) Isopropyl Acetate	8.694	43	294456	29.231	ug/l #	82
51) 4-Methyl-2-Pentanone	10.453	43	294262	49.111	ug/l	91

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

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