

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078987.D
 Acq On : 22 Aug 2023 12:22
 Operator : JC\MD
 Sample : PB154968ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#02

Quant Time: Aug 23 03:00:20 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	180658	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	334895	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	336388	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	99419	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	143256	50.959	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 101.920%		
35) Dibromofluoromethane	8.171	113	117738	52.227	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 104.460%		
50) Toluene-d8	10.571	98	432921	51.720	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.440%		
62) 4-Bromofluorobenzene	12.853	95	133520	48.801	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 97.600%		
Target Compounds						
						Qvalue
16) Acetone	4.442	43	25684	22.143	ug/l	88
20) Methylene Chloride	5.283	84	11304	4.573	ug/l #	82
25) 2-Butanone	7.494	43	10704	5.756	ug/l #	52
37) Ethyl Acetate	7.571	43	5544	1.461	ug/l #	79
43) Isopropyl Acetate	8.694	43	159301	23.754	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	39350	9.893	ug/l #	83

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

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