

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078838.D
 Acq On : 10 Aug 2023 15:32
 Operator : JC\MD
 Sample : PB154726ZHE#09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#09

Quant Time: Aug 10 23:52:48 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	176793	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	340414	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	314921	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	82647	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	141576	53.417	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.840%	
35) Dibromofluoromethane	8.177	113	115112	50.074	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.140%	
50) Toluene-d8	10.571	98	436514	51.599	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 103.200%	
62) 4-Bromofluorobenzene	12.859	95	113507	44.895	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 89.780%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	67141	65.840	ug/l #	83
20) Methylene Chloride	5.289	84	13628	5.396	ug/l #	91
25) 2-Butanone	7.494	43	32767	18.975	ug/l #	80
37) Ethyl Acetate	7.571	43	33098	8.728	ug/l	93
43) Isopropyl Acetate	8.700	43	197463	30.886	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	111589	28.496	ug/l #	85

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

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