

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078991.D
 Acq On : 22 Aug 2023 13:59
 Operator : JC\MD
 Sample : PB154968ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#06

Quant Time: Aug 23 03:02:09 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	176568	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	338571	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	337304	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	105576	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	145600	52.993	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 105.980%	
35) Dibromofluoromethane	8.171	113	117000	51.336	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.680%	
50) Toluene-d8	10.571	98	440563	52.061	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.120%	
62) 4-Bromofluorobenzene	12.853	95	136257	49.261	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 98.520%	
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	28601	25.228	ug/l	89
20) Methylene Chloride	5.289	84	8293	3.433	ug/l #	81
25) 2-Butanone	7.500	43	15430	8.489	ug/l #	71
43) Isopropyl Acetate	8.694	43	195465	28.831	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	65030	16.172	ug/l #	81

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

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