

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082223\
 Data File : VN078990.D
 Acq On : 22 Aug 2023 13:35
 Operator : JC\MD
 Sample : PB154968ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154968ZHE#05

Quant Time: Aug 23 03:01:41 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	175667	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	343370	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	348237	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	108629	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	143419	52.467	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.940%	
35) Dibromofluoromethane	8.171	113	121032	52.363	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.720%	
50) Toluene-d8	10.571	98	444751	51.821	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 103.640%	
62) 4-Bromofluorobenzene	12.853	95	139958	49.892	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 99.780%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	29132	25.829	ug/l #	78
20) Methylene Chloride	5.283	84	8220	3.420	ug/l #	78
25) 2-Butanone	7.494	43	17167	9.493	ug/l #	70
43) Isopropyl Acetate	8.700	43	196320	28.552	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	65714	16.114	ug/l #	80

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

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