

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082923\
 Data File : VN079103.D
 Acq On : 29 Aug 2023 18:56
 Operator : JC\MD
 Sample : PB155153ZHE#12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155153ZHE#12

Quant Time: Aug 29 23:32:13 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	137585	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	275513	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	275989	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	82668	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	118022	55.126	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 110.260%	
35) Dibromofluoromethane	8.177	113	94694	51.059	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.120%	
50) Toluene-d8	10.571	98	365214	53.035	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 106.060%	
62) 4-Bromofluorobenzene	12.853	95	103805	46.118	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.240%	
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	25194	28.520	ug/l	89
20) Methylene Chloride	5.283	84	16060	8.531	ug/l #	87
25) 2-Butanone	7.500	43	7460	5.267	ug/l #	61
43) Isopropyl Acetate	8.694	43	210218	38.103	ug/l #	71
51) 4-Methyl-2-Pentanone	10.447	43	27763	8.485	ug/l #	71

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

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