

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082923\
 Data File : VN079100.D
 Acq On : 29 Aug 2023 17:43
 Operator : JC\MD
 Sample : PB155153ZHE#09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155153ZHE#09

Quant Time: Aug 29 23:31:34 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	141950	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	275226	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	276337	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	84912	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	117752	53.309	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.620%	
35) Dibromofluoromethane	8.171	113	95389	51.487	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 102.980%	
50) Toluene-d8	10.571	98	360547	52.412	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.820%	
62) 4-Bromofluorobenzene	12.853	95	107072	47.619	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 95.240%	
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	23504	25.789	ug/l	89
20) Methylene Chloride	5.277	84	16622	8.558	ug/l #	76
43) Isopropyl Acetate	8.694	43	199300	36.162	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	28095	8.595	ug/l #	73

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

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