

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082923\
 Data File : VN079102.D
 Acq On : 29 Aug 2023 18:31
 Operator : JC\MD
 Sample : PB155153ZHE#11
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155153ZHE#11

Quant Time: Aug 29 23:32:00 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	8.229	168	136666	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	274953	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	277765	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84886	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	116083	54.585	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.180%	
35) Dibromofluoromethane	8.171	113	93828	50.695	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.380%	
50) Toluene-d8	10.570	98	359641	52.332	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.660%	
62) 4-Bromofluorobenzene	12.853	95	106543	47.431	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 94.860%	
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	24228	27.611	ug/l	94
20) Methylene Chloride	5.277	84	17696	9.463	ug/l #	89
43) Isopropyl Acetate	8.694	43	205175	37.265	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	27008	8.271	ug/l #	71

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

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