

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN082923\
 Data File : VN079098.D
 Acq On : 29 Aug 2023 16:54
 Operator : JC\MD
 Sample : PB155153ZHE#07
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155153ZHE#07

Quant Time: Aug 29 23:31:06 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	135215	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	273344	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	276537	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	82380	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	113499	53.943	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.880%	
35) Dibromofluoromethane	8.177	113	92894	50.486	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.980%	
50) Toluene-d8	10.571	98	355926	52.096	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 104.200%	
62) 4-Bromofluorobenzene	12.853	95	102761	46.016	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 92.040%	
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	25721	29.627	ug/l	97
20) Methylene Chloride	5.277	84	22593	12.212	ug/l #	82
25) 2-Butanone	7.500	43	7779	5.589	ug/l #	57
43) Isopropyl Acetate	8.694	43	189977	34.708	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	31583	9.729	ug/l #	74

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

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