

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090223\
 Data File : VN079143.D
 Acq On : 02 Sep 2023 15:11
 Operator : JC\MD
 Sample : PB155301ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB155301ZHE#05

Quant Time: Sep 04 01:04:31 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	210027	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	411274	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	411521	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	124376	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	158625	48.536	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	97.080%
35) Dibromofluoromethane	8.171	113	139375	50.343	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.680%
50) Toluene-d8	10.571	98	533269	51.876	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.760%
62) 4-Bromofluorobenzene	12.853	95	162385	48.329	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.660%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	28175	20.893	ug/l	92
20) Methylene Chloride	5.283	84	19681	6.849	ug/l #	81
43) Isopropyl Acetate	8.694	43	217178	26.371	ug/l #	77
51) 4-Methyl-2-Pentanone	10.453	43	29480	6.035	ug/l #	82

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

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