

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078949.D
 Acq On : 18 Aug 2023 18:00
 Operator : JC\MD
 Sample : PB154902ZHE#06
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154902ZHE#06

Quant Time: Aug 18 23:46:02 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	146052	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	281855	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	271044	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	75509	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	116988	51.475	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	102.960%
35) Dibromofluoromethane	8.171	113	97894	51.596	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.200%
50) Toluene-d8	10.571	98	359601	51.045	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	102.080%
62) 4-Bromofluorobenzene	12.853	95	101994	44.294	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	88.580%
Target Compounds						
						Qvalue
16) Acetone	4.442	43	29226	31.166	ug/l	92
20) Methylene Chloride	5.289	84	10716	5.362	ug/l	89
25) 2-Butanone	7.500	43	12017	7.993	ug/l #	77
43) Isopropyl Acetate	8.694	43	224971	39.860	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	45983	13.737	ug/l #	81

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

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