

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081823\
 Data File : VN078945.D
 Acq On : 18 Aug 2023 16:23
 Operator : JC\MD
 Sample : PB154902ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154902ZHE#02

Quant Time: Aug 18 23:44:03 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	151217	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	278716	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	275658	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84117	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	117133	49.779	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.560%
35) Dibromofluoromethane	8.171	113	94830	50.544	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.080%
50) Toluene-d8	10.571	98	360983	51.818	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	103.640%
62) 4-Bromofluorobenzene	12.853	95	107913	47.392	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	94.780%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	31279	32.216	ug/l #	84
20) Methylene Chloride	5.283	84	17527	8.471	ug/l	86
25) 2-Butanone	7.494	43	9372	6.021	ug/l #	67
43) Isopropyl Acetate	8.694	43	170714	30.587	ug/l #	79
51) 4-Methyl-2-Pentanone	10.447	43	41820	12.634	ug/l #	79

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

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