

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078846.D
 Acq On : 10 Aug 2023 18:45
 Operator : JC\MD
 Sample : PB154726ZHE#17
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#17

Quant Time: Aug 10 23:55:59 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080723W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 08 05:03:52 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.236	168	169565	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	342995	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	310511	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	89744	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	142319	55.986	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 111.980%		
35) Dibromofluoromethane	8.177	113	114685	49.512	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 99.020%		
50) Toluene-d8	10.571	98	432078	50.690	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 101.380%		
62) 4-Bromofluorobenzene	12.859	95	120558	47.325	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 94.640%		
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	67543	69.057	ug/l #	85
20) Methylene Chloride	5.295	84	13885	5.732	ug/l #	78
25) 2-Butanone	7.494	43	32359	19.538	ug/l #	75
37) Ethyl Acetate	7.577	43	32868	8.603	ug/l #	92
43) Isopropyl Acetate	8.694	43	243308	37.771	ug/l #	69
51) 4-Methyl-2-Pentanone	10.453	43	115901	29.374	ug/l #	83

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

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