

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078845.D
 Acq On : 10 Aug 2023 18:21
 Operator : JC\MD
 Sample : PB154726ZHE#16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#16

Quant Time: Aug 10 23:55:35 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	8.235	168	169754	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	332884	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	312122	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	90438	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	135555	53.266	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 106.540%		
35) Dibromofluoromethane	8.177	113	109907	48.891	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.780%		
50) Toluene-d8	10.571	98	431794	52.196	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 104.400%		
62) 4-Bromofluorobenzene	12.859	95	117663	47.591	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.180%		
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	67381	68.815	ug/l #	86
20) Methylene Chloride	5.289	84	15697	6.473	ug/l #	78
25) 2-Butanone	7.494	43	31777	19.165	ug/l #	86
37) Ethyl Acetate	7.565	43	37112	10.008	ug/l #	88
43) Isopropyl Acetate	8.694	43	239954	38.382	ug/l #	70
51) 4-Methyl-2-Pentanone	10.447	43	115739	30.224	ug/l #	81

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

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