

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078834.D
 Acq On : 10 Aug 2023 13:55
 Operator : JC\MD
 Sample : PB154726ZHE#05
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#05

Quant Time: Aug 10 23:51:23 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	167872	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	330186	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	304917	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	78140	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	132470	52.637	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 105.280%		
35) Dibromofluoromethane	8.177	113	112041	50.247	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.500%		
50) Toluene-d8	10.571	98	421793	51.403	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 102.800%		
62) 4-Bromofluorobenzene	12.853	95	115200	46.976	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 93.960%		
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	64433	66.542	ug/l	97
20) Methylene Chloride	5.300	84	11867	4.948	ug/l #	83
25) 2-Butanone	7.500	43	31004	18.908	ug/l #	74
37) Ethyl Acetate	7.577	43	34181	9.293	ug/l #	88
43) Isopropyl Acetate	8.700	43	171482	27.653	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	112964	29.741	ug/l #	82

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

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