

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078842.D
 Acq On : 10 Aug 2023 17:09
 Operator : JC\MD
 Sample : PB154726ZHE#13
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#13

Quant Time: Aug 10 23:54:22 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	170199	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	339530	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	312876	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84337	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	138426	54.252	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 108.500%		
35) Dibromofluoromethane	8.177	113	111608	48.676	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 97.360%		
50) Toluene-d8	10.571	98	433012	51.318	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 102.640%		
62) 4-Bromofluorobenzene	12.853	95	116622	46.247	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 92.500%		
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	68445	69.719	ug/l	94
20) Methylene Chloride	5.300	84	14860	6.112	ug/l #	79
25) 2-Butanone	7.500	43	33866	20.371	ug/l #	73
37) Ethyl Acetate	7.577	43	36051	9.532	ug/l #	91
43) Isopropyl Acetate	8.700	43	238888	37.463	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	116055	29.714	ug/l #	83

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

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