

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081023\
 Data File : VN078844.D
 Acq On : 10 Aug 2023 17:57
 Operator : JC\MD
 Sample : PB154726ZHE#15
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154726ZHE#15

Quant Time: Aug 10 23:55:07 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	166409	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	333483	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	304732	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84675	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	139552	55.939	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 111.880%			
35) Dibromofluoromethane	8.177	113	110491	49.062	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 98.120%			
50) Toluene-d8	10.571	98	421739	50.888	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 101.780%			
62) 4-Bromofluorobenzene	12.853	95	116614	47.082	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 94.160%			
Target Compounds					Qvalue	
16) Acetone	4.436	43	68199	71.050	ug/l	90
20) Methylene Chloride	5.289	84	15099	6.351	ug/l #	71
25) 2-Butanone	7.494	43	33244	20.453	ug/l #	80
37) Ethyl Acetate	7.571	43	36972	9.953	ug/l #	92
43) Isopropyl Acetate	8.700	43	233228	37.239	ug/l #	70
51) 4-Methyl-2-Pentanone	10.453	43	117807	30.709	ug/l #	82

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

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