

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078901.D
 Acq On : 16 Aug 2023 14:28
 Operator : JC\MD
 Sample : PB154855ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154855ZHE#04

Quant Time: Aug 17 01:16:32 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	150701	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	304646	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	299551	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	93428	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	130551	55.671	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 111.340%		
35) Dibromofluoromethane	8.171	113	102752	50.105	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.220%		
50) Toluene-d8	10.571	98	392612	51.561	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.120%		
62) 4-Bromofluorobenzene	12.853	95	121249	48.716	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 97.440%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	28087	29.028	ug/l #	75
20) Methylene Chloride	5.283	84	13670	6.629	ug/l #	73
25) 2-Butanone	7.500	43	8172	5.268	ug/l #	78
37) Ethyl Acetate	7.571	43	3708	1.074	ug/l #	81
43) Isopropyl Acetate	8.694	43	220428	36.133	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	35456	9.799	ug/l #	82

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

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