

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078898.D
 Acq On : 16 Aug 2023 13:16
 Operator : JC\MD
 Sample : PB154855ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154855ZHE#01

Quant Time: Aug 17 01:15:23 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	155650	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	299075	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	294985	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	91756	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	127229	52.530	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 105.060%		
35) Dibromofluoromethane	8.171	113	100987	50.162	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 100.320%		
50) Toluene-d8	10.571	98	397715	53.204	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 106.400%		
62) 4-Bromofluorobenzene	12.853	95	116996	47.883	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 95.760%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	88480	88.535	ug/l	88
20) Methylene Chloride	5.277	84	25530	11.987	ug/l #	89
25) 2-Butanone	7.500	43	40266	25.131	ug/l #	81
37) Ethyl Acetate	7.565	43	43301	12.779	ug/l #	90
43) Isopropyl Acetate	8.694	43	174468	29.132	ug/l #	78
51) 4-Methyl-2-Pentanone	10.447	43	138783	39.072	ug/l #	83

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

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