

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078899.D
 Acq On : 16 Aug 2023 13:40
 Operator : JC\MD
 Sample : PB154855ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154855ZHE#02

Quant Time: Aug 17 01:15:45 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	148917	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	288702	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	282270	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	84686	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	125934	54.346	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 108.700%		
35) Dibromofluoromethane	8.171	113	100055	51.485	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 102.960%		
50) Toluene-d8	10.571	98	384787	53.324	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 106.640%		
62) 4-Bromofluorobenzene	12.853	95	111977	47.476	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 94.960%		
Target Compounds						
					Qvalue	
16) Acetone	4.430	43	77679	81.242	ug/l	89
20) Methylene Chloride	5.283	84	23287	11.429	ug/l #	78
25) 2-Butanone	7.494	43	37198	24.266	ug/l #	81
37) Ethyl Acetate	7.571	43	40709	12.446	ug/l #	88
43) Isopropyl Acetate	8.694	43	191312	33.092	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	127214	37.101	ug/l #	82

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

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