

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081623\
 Data File : VN078897.D
 Acq On : 16 Aug 2023 12:52
 Operator : JC\MD
 Sample : PB154855TB
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154855TB

Quant Time: Aug 17 01:14:56 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	163119	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	306102	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	303749	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	93121	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	129828	51.148	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 102.300%		
35) Dibromofluoromethane	8.171	113	108763	52.784	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 105.560%		
50) Toluene-d8	10.571	98	412020	53.853	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 107.700%		
62) 4-Bromofluorobenzene	12.853	95	125075	50.015	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 100.020%		
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	83605	79.827	ug/l #	85
20) Methylene Chloride	5.289	84	16588	7.432	ug/l #	86
25) 2-Butanone	7.494	43	41559	24.750	ug/l #	80
37) Ethyl Acetate	7.571	43	42855	12.357	ug/l #	89
43) Isopropyl Acetate	8.694	43	175678	28.661	ug/l #	78
51) 4-Methyl-2-Pentanone	10.453	43	142177	39.108	ug/l #	82

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

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