

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078665.D
 Acq On : 27 Jul 2023 17:20
 Operator : JC\MD
 Sample : PB154443ZHE#02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154443ZHE#02

Quant Time: Jul 28 01:09:35 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N071023W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 10 16:13:53 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	423742	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	772665	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	751382	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	280815	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	274679	48.353	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	=	96.700%	
35) Dibromofluoromethane	8.177	113	275165	52.461	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	104.920%	
50) Toluene-d8	10.571	98	955807	48.238	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	=	96.480%	
62) 4-Bromofluorobenzene	12.853	95	327926	53.730	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	107.460%	
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	93881	53.054	ug/l	90
20) Methylene Chloride	5.289	84	27082	4.216	ug/l #	94
25) 2-Butanone	7.506	43	51940	18.647	ug/l	96
37) Ethyl Acetate	7.577	43	14073	2.229	ug/l #	95
43) Isopropyl Acetate	8.700	43	350093	33.682	ug/l #	83
51) 4-Methyl-2-Pentanone	10.453	43	174075	28.156	ug/l	91

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

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