

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078564.D
 Acq On : 18 Jul 2023 16:27
 Operator : JC\MD
 Sample : PB154251ZHE#01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154251ZHE#01

Quant Time: Jul 19 01:52:37 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	407702	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	763834	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	736206	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	272262	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	272487	49.854	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	99.700%
35) Dibromofluoromethane	8.171	113	276025	53.234	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	106.460%
50) Toluene-d8	10.577	98	1046839	53.443	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	106.880%
62) 4-Bromofluorobenzene	12.853	95	319993	53.036	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	106.080%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	119522	70.201	ug/l	99
20) Methylene Chloride	5.283	84	25634	4.127	ug/l #	91
25) 2-Butanone	7.494	43	87396	32.610	ug/l #	87
37) Ethyl Acetate	7.577	43	22136	3.547	ug/l	98
43) Isopropyl Acetate	8.700	43	374596	36.456	ug/l #	76
51) 4-Methyl-2-Pentanone	10.453	43	297567	48.687	ug/l	90

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

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