

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072723\
 Data File : VN078667.D
 Acq On : 27 Jul 2023 18:09
 Operator : JC\MD
 Sample : PB154443ZHE#04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB154443ZHE#04

Quant Time: Jul 28 01:10:24 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	400672	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	729576	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	695957	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	255475	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.589	65	258127	48.056	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	96.120%
35) Dibromofluoromethane	8.177	113	261121	52.724	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	105.440%
50) Toluene-d8	10.571	98	891630	47.657	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	95.320%
62) 4-Bromofluorobenzene	12.853	95	298309	51.764	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	103.520%
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	94259	56.334	ug/l	100
20) Methylene Chloride	5.289	84	27592	4.640	ug/l	99
25) 2-Butanone	7.500	43	54748	20.786	ug/l	91
37) Ethyl Acetate	7.571	43	17217	2.888	ug/l #	95
43) Isopropyl Acetate	8.700	43	516695	52.646	ug/l #	73
51) 4-Methyl-2-Pentanone	10.453	43	176974	30.315	ug/l	90

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

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