

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062123\
 Data File : VN078339.D
 Acq On : 21 Jun 2023 21:59
 Operator : JC\MD
 Sample : 03357-10
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-SCAN-02-G

Quant Time: Jun 22 04:57:28 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060623W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 07 05:05:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.236	168	472731	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	906119	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	826698	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	259522	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	340373	54.002	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 108.000%			
35) Dibromofluoromethane	8.177	113	309538	54.027	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 108.060%			
50) Toluene-d8	10.577	98	1085301	49.138	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.280%			
62) 4-Bromofluorobenzene	12.859	95	339832	49.493	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 98.980%			
Target Compounds						
					Qvalue	
16) Acetone	4.448	43	172879	95.291	ug/l	98
20) Methylene Chloride	5.289	84	13404	2.236	ug/l #	86
25) 2-Butanone	7.494	43	49129	17.610	ug/l #	79
37) Ethyl Acetate	7.577	43	26613	4.145	ug/l	96
43) Isopropyl Acetate	8.694	43	215187	17.527	ug/l #	72
51) 4-Methyl-2-Pentanone	10.453	43	321794	49.772	ug/l	91

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

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