

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071423\
 Data File : VN078519.D
 Acq On : 14 Jul 2023 13:58
 Operator : JC\MD
 Sample : 03580-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WASTE-1

Quant Time: Jul 15 04:01:19 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

Manual Integrations
 APPROVED

Reviewed By : John Carlone 07/17/2023
 Supervised By : Mahesh Dadoda 07/17/2023

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

Internal Standards						
1) Pentafluorobenzene	8.235	168	400785	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	748200	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	724846	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	257533	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	274931	51.169	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 102.340%			
35) Dibromofluoromethane	8.171	113	266052	52.383	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 104.760%			
50) Toluene-d8	10.576	98	896587	46.729	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 93.460%			
62) 4-Bromofluorobenzene	12.859	95	305859	51.753	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 103.500%			
Target Compounds						
					Qvalue	
16) Acetone	4.441	43	139935	83.609	ug/l	89
20) Methylene Chloride	5.294	84	25981m	4.294	ug/l	
25) 2-Butanone	7.500	43	79485	30.170	ug/l #	83
37) Ethyl Acetate	7.582	43	22363	3.658	ug/l #	78
43) Isopropyl Acetate	8.700	43	236826	23.530	ug/l #	84
51) 4-Methyl-2-Pentanone	10.453	43	263258	43.973	ug/l	91

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

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