

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071823\
 Data File : VN078563.D
 Acq On : 18 Jul 2023 16:03
 Operator : JC\MD
 Sample : 03654-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-12016

Quant Time: Jul 19 01:52:05 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	494971	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	901922	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	896139	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	312554	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	308422	46.480	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	92.960%
35) Dibromofluoromethane	8.171	113	333904	54.537	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	109.080%
50) Toluene-d8	10.577	98	982032	42.459	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	84.920%#
62) 4-Bromofluorobenzene	12.853	95	362668	50.906	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.820%
Target Compounds						
					Qvalue	
16) Acetone	4.442	43	66599	32.220	ug/l #	85
20) Methylene Chloride	5.289	84	19475	2.114	ug/l	91
25) 2-Butanone	7.500	43	50150	15.413	ug/l	99
37) Ethyl Acetate	7.577	43	15119	2.052	ug/l #	91
43) Isopropyl Acetate	8.700	43	545290	44.943	ug/l #	74
51) 4-Methyl-2-Pentanone	10.453	43	160936	22.300	ug/l	90

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

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