

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062223\
 Data File : VN078358.D
 Acq On : 22 Jun 2023 18:28
 Operator : JC\MD
 Sample : 03361-16
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TCLP-VOC-SB-23-GRAB-5

Quant Time: Jun 23 08:03:57 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	8.235	168	467755	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	873905	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	813702	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	263717	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	326452	52.345	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 104.680%	
35) Dibromofluoromethane	8.177	113	304900	55.180	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 110.360%	
50) Toluene-d8	10.571	98	1039660	48.806	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 97.620%	
62) 4-Bromofluorobenzene	12.853	95	335216	50.620	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 101.240%	
Target Compounds						
					Qvalue	
16) Acetone	4.447	43	65561	36.522	ug/l	98
20) Methylene Chloride	5.300	84	24114	4.065	ug/l	93
25) 2-Butanone	7.506	43	18122	6.565	ug/l #	81
37) Ethyl Acetate	7.571	43	51777	8.362	ug/l	98
43) Isopropyl Acetate	8.694	43	279540	23.607	ug/l #	82
51) 4-Methyl-2-Pentanone	10.453	43	143935	23.083	ug/l	88

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

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