

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN060723\
 Data File : VN078109.D
 Acq On : 07 Jun 2023 17:27
 Operator : JC\MD
 Sample : 03067-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 DT-GW-1

Quant Time: Jun 08 05:29:20 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	355427	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.112	114	652218	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	558224	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	169119	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	243152	51.309	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 102.620%	
35) Dibromofluoromethane	8.171	113	212607	51.555	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 103.100%	
50) Toluene-d8	10.570	98	804032	50.574	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 101.140%	
62) 4-Bromofluorobenzene	12.853	95	224361	45.396	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 90.800%	
Target Compounds						
						Qvalue
5) Bromomethane	2.989	94	1627	0.409	ug/l	99
16) Acetone	4.447	43	14540	10.660	ug/l	92
18) Methyl Acetate	5.041	43	5833	0.929	ug/l #	71
52) Toluene	10.635	92	26675	2.392	ug/l	93
68) m/p-Xylenes	12.070	106	3895	0.495	ug/l	85
69) o-Xylene	12.400	106	1723	0.223	ug/l	89
84) 1,2,4-Trimethylbenzene	13.494	105	3990	0.318	ug/l	99

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

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