

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061923\
 Data File : VN078289.D
 Acq On : 19 Jun 2023 19:26
 Operator : JC\MD
 Sample : 03320-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-12

Quant Time: Jun 20 01:33:40 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	459771	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	867593	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.870	117	813327	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.800	152	269197	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.588	65	332164	54.185	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 108.380%		
35) Dibromofluoromethane	8.177	113	294530	53.691	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 107.380%		
50) Toluene-d8	10.576	98	1032653	48.830	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 97.660%		
62) 4-Bromofluorobenzene	12.853	95	345738	52.589	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 105.180%		
Target Compounds						
					Qvalue	
16) Acetone	4.453	43	25975	14.721	ug/l	94
20) Methylene Chloride	5.288	84	15828	2.715	ug/l #	86
37) Ethyl Acetate	7.571	43	63266	10.292	ug/l	96
43) Isopropyl Acetate	8.700	43	413642	35.186	ug/l #	76

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061923\
Data File : VN078289.D
Acq On : 19 Jun 2023 19:26
Operator : JC\MD
Sample : 03320-08
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-12

Quant Time: Jun 20 01:33:40 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

