

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010324\
 Data File : VN080643.D
 Acq On : 03 Jan 2024 16:31
 Operator : JC\MD
 Sample : P1003-02
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 72-11978

Quant Time: Jan 04 02:01:46 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122723W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 28 03:07:01 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.224	168	285909	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	561851	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	627505	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	274656	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	185694	50.633	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 101.260%		
35) Dibromofluoromethane	8.165	113	173278	47.685	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 95.380%		
50) Toluene-d8	10.565	98	653599	51.515	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 103.040%		
62) 4-Bromofluorobenzene	12.853	95	258157	60.459	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 120.920%		
Target Compounds						
					Qvalue	
16) Acetone	4.418	43	27329	27.413	ug/l	99
20) Methylene Chloride	5.277	84	21205	6.666	ug/l	90
25) 2-Butanone	7.488	43	24004	12.883	ug/l	92
43) Isopropyl Acetate	8.688	43	97583	14.267	ug/l #	87

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010324\
Data File : VN080643.D
Acq On : 03 Jan 2024 16:31
Operator : JC\MD
Sample : P1003-02
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
72-11978

Quant Time: Jan 04 02:01:46 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122723W.M
Quant Title : SW846 8260
QLast Update : Thu Dec 28 03:07:01 2023
Response via : Initial Calibration

