

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102021\
 Data File : VN069060.D
 Acq On : 20 Oct 2021 14:30
 Operator : JC/MD
 Sample : M4259-04
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-1-BS

Quant Time: Oct 20 23:14:19 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 13 05:44:40 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.088	168	339789	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.971	114	594798	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.747	117	569798	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.678	152	212514	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.440	65	222957	49.767	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	99.540%		
35) Dibromofluoromethane	8.024	113	175783	50.796	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	101.600%		
50) Toluene-d8	10.443	98	743393	53.689	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	107.380%		
62) 4-Bromofluorobenzene	12.734	95	258856	49.551	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	99.100%		
Target Compounds						
					Qvalue	
16) Acetone	4.299	43	14605	8.736	ug/l	100
18) Methyl Acetate	4.865	43	3119	0.492	ug/l #	81
20) Methylene Chloride	5.103	84	20651	5.367	ug/l	94
43) Isopropyl Acetate	8.563	43	35549	3.840	ug/l #	91

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN102021\
Data File : VN069060.D
Acq On : 20 Oct 2021 14:30
Operator : JC/MD
Sample : M4259-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
WC-1-BS

Quant Time: Oct 20 23:14:19 2021
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N101221W.M
Quant Title : SW846 8260
QLast Update : Wed Oct 13 05:44:40 2021
Response via : Initial Calibration

